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CFL3D Version 6.4—General Usage and Aeroelastic Analysis

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Available from:



Abstract

This document contains the course notes on the computational fluid dynamics code CFL3D version 6.4. It is intended to provide from basic to advanced users the information necessary to successfully use the code for a broad range of cases. Much of the course covers capability that has been a part of previous versions of the code, with material compiled from a CFL3D v5.0 manual and from the CFL3D v6 web site prior to the current release. This part of the material is presented to users of the code not familiar with computational fluid dynamics. There is new capability in CFL3D version 6.4 presented here that has not previously been published. There are also outdated features no longer used or recommended in recent releases of the code. The information offered here supersedes earlier manuals and updates outdated usage. Where current usage supersedes older versions, notation of that is made. These course notes also provides hints for usage, code installation and examples not found elsewhere.

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Introduction



CFL3D is a Reynolds-averaged Navier-Stokes flow solver for structured grids. The original version, developed in the 1980's was given its name to denote its origin in the Computational Fluids Laboratory. CFL3D solves the time-dependent conservation law form of the equations using a semi-discrete finite-volume approach with upwind-biasing of the convective and pressure terms and central differencing of the shear stress and heat transfer terms. Numerous turbulence models are provided. Grids must be supplied external to the code.

The present document is an outgrowth of a course that was presented on the computational fluid dynamics code CFL3D version 6.4. Publication of this material in the present form makes it available to many more users of the code. This document should provide the information necessary to successfully use the code for a broad range of cases. The target audience ranges from basic to advanced users. New users should find useful the discussion of general features of the code and the many options that are available, code set up, creation of grids and input for steady and unsteady computations. New features that are available in CFL3D version 6.4 will also be discussed. There is a lengthy discussion of issues related to unsteady computations, moving and deforming meshes, aeroelastic simulations and parallel computing using the message passing interface (MPI). Within these discussions there are detailed instructions on input parameters, their use within the code, as well as illustrative examples.

Much of the course covers capability that has been a part of previous versions of the code, with material compiled from a CFL3D v5.0 manual and from the CFL3D v6 web site prior to the current release. This part of the material is presented to users of the code not familiar with computational fluid dynamics. There is also new capability in CFL3D v6.4 that has not previously been published. This course intends to acquaint users with this new capability. There are also outdated features no longer used or recommended in recent releases of the code. The information offered here supersedes earlier manuals and updates outdated usage. Where current usage supersedes older versions, notation of that is made. This document also provides hints for usage and code installation not found elsewhere.

Introduction



There is much information in the CFL3D v5.0 manual that is not presented in these notes. The use of patched, overset or embedded grids is not discussed here. Since the intention is to provide users a practical guide on code usage, there is very little discussion of the fluid dynamics equations and computational method used. This information is available in the CFL3D v5.0 manual.

The attempt is to organize this material in an intuitive way. Topics are presented in the order they would be encountered in the process of building up a real test case. The ordering of the information reflects the course instructor's own learning experience with CFL3D. Others may order the material differently. This course is not comprehensive. Because of the vast number of ways in which CFL3D can be used there are many input options that are not discussed and none are discussed in complete detail. Those that are discussed are the more commonly used features. By the end of the course the reader should be able to perform a number of different analyses with the code. If the reader is interested in more detail also consult the CFL3D v6 web page and the CFL3D v5.0 user's manual. These references are listed at the back of this document.

What's New in CFL3D v6.4



There are new capabilities in CFL3D v6.4 presented in this document. They are:

- New mesh deformation scheme with more options available.
- New aeroelastic analyses not available in previous versions of CFL3D
- 2nd order time accuracy in turbulence modeling (default)
- New keywords are available
 - 1st time accurate turbulence modeling (default is 2nd order)
 - New options in turbulence modeling
 - Full Navier-Stokes terms available
 - Option to exercise mesh deformation without full flow solver
 - Calculation of CFL number can be modified for axisymmetric cases to increase convergence rate
- Changes in the input for prescribed modal motion



Major features

- Euler
- Laminar thin-layer Navier-Stokes
- Reynolds-Averaged thin-layer Navier-Stokes (RANS)
- Structured grid
- Single or multi-block
- Dynamic memory
- Parallel (MPI) capability
- Moving grid and mesh deformation capability
- CGNS (CFD General Notation System) capability for CFD output

Discretization and numerical method

- Conservation law form of the Euler or RANS equations
- Spatial discretization is semi-discrete finite-volume approach
- Upwind-Biasing is used for the convective and pressure terms
- Solves either the steady or unsteady form of the equations
- Time advancement is implicit with dual time stepping and sub-iterations



- Discretization and numerical method (...continued)
 - Approximate-Factorized (AF) numerical scheme
 - Explicit block boundary conditions
 - Multigrid
 - Grid sequencing
- Block structures
 - 1-1 blocking (preferred)
 - Patching
 - Overlapping
 - Embedding
 - Sliding patched zone interfaces
 - Grids must have been created prior to execution of CFL3D



- Turbulence models for RANS computation
 - 0-equation models: Baldwin-Lomax, Baldwin-Lomax with Degani-Schiff modification
 - 1-equation models: Baldwin-Barth, Spalart-Almaras, including Detached Eddy Simulation (DES)
 - 2-equation models: Wilcox k- ω model, Menter's k- ω Shear Stress Transport (SST) model, Abid k- ω model, k- ω and k- ϵ Explicit Algebraic Stress Models (EASM), k-enstrophy model
- Computing modes
 - Sequential or single processor (single or multiple blocks)
 - Parallel processing
 - Message Passing Interface (MPI)
 - Requires multi-block structure
 - May be run on distributed memory machines. (PC clusters or parallel supercomputer)



- Computing modes (...continued)
 - Complex computation
 - Allows computation of sensitivity derivatives due to static and dynamic variables (e.g. $dC_1/d\alpha$)
 - Requires compiling of the complex executable for static and dynamic sensitivity calculations
 - Dynamic sensitivity calculations require additional keyword input
- Code developers and points of contact:
 - Many developers have contributed to CFL3D
 - Most recent primary NASA LaRC developers (POC's) are:
 - Dr. Robert T. Biedron (757-864-2156, r.t.biedron@larc.nasa.gov) general flow solver, multiblock, MPI
 - Dr. Christopher Rumsey (757-864-2165,c.l.rumsey@larc.nasa.gov) turbulence models
 - Dr. Bob Bartels (757-864-2813, r.e.bartels@larc.nasa.gov) aeroelastic modules and deforming mesh



- Online and printable documentation: http://cfl3d.larc.nasa.gov/Cfl3dv6/cfl3dv6.html
 - Recommend printing the Version 5.0 manual for reference (found as a link at the web site above)
- Acquiring the code:
 - Version 6 is currently available for general distribution to U.S. citizens within the United States.
 The code cannot be released outside of the United States. If you would like a copy of the code, please follow the request procedure below:
 - Send <u>e-mail</u> or FAX (757-864-8816) to one of the POC's requesting CFL3D Version 6, along with a brief description of your planned usage of the code, your phone number, and FAX number.
 - Your request will be forwarded internally to a NASA Software Releasing Authority (SRA). The SRA will determine whether or not the code may be released to you; if so, the SRA will e-mail or FAX a Usage Agreement to you to fill out, sign and return to the SRA.



- After the SRA has granted permission, the code will be provided to you
 electronically. In addition, you will be added to the Version 6 user list, and will
 receive any updates and/or corrections that occur.
- Note: even if you are a registered Version 5 user you must still follow the formal request procedure for Version 6.
- Conditions of use:
 - Do not distribute any part of the code outside of your working group
 - Report any bugs you may find
 - CFL3D is restricted to use within the United States
 - Abide by any additional conditions in the usage agreement



 To install CFL3Dv6 on a particular machine, you must have the following file:

cfl3dv6.tar.DATE.gz (tarred and gzipped version 6 package)

Note: DATE indicates the release date in the form MMM_DD_YYYY. For example, cfl3dv6.tar.Sep_12_2003 indicates the code as of September 12, 2003.

 Make sure that: ./ is in your path; if not, you will have to explicitly prepend ./ to all the commands below

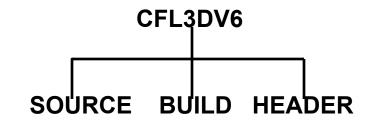
Type:

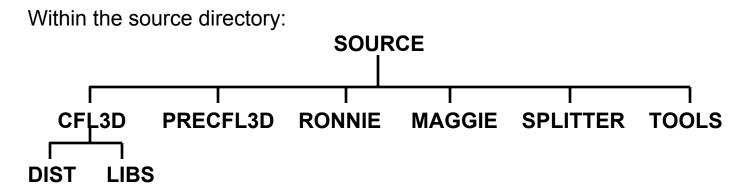
gunzip cfl3dv6.tar.DATE.gz

tar -xvf cfl3dv6.tar.DATE



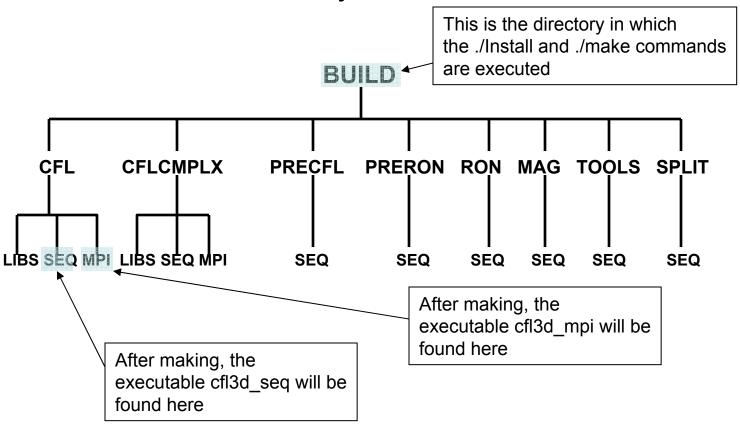
You should end up with the following directory structure:







Within the build directory:





– In the subdirectory build, type:

Install [options] or ./Install [options]

Where [options] may be blank or one or more of the following:

- -no opt
- create executables with little optimization but fast compilation
- -single
- create single precision executables
- -noredirect
- · disallow redirected input file; needed only for SP2 and sometimes on Linux with MPI
- -mpichdir=dir1
- use MPICH on a workstation cluster; dir1 is the directory where mpich is located not used on MPP machines
- -linux compiler flags=flag
- sets up to compile using special compiler flags for use on Linux operating systems only; flag is currently Intel, PG, Lahey, or Alpha (Intel is currently the default) Example: To use the Portland Group compiler MUST install with: ./Install -linux_compiler_flags=PG
- -help
- print out the Install options



- Note: the directory paths for either the mpichdir or cgnsdir options should be either absolute paths or paths relative to the installation directory; the use of ~ to denote a home directory is not allowed.
- If -no_opt is not specified, various compiler optimization levels are used to speed execution but results in slower compilation.
- If -mpichdir=dir1 is not used, then it is assumed "native" MPI is available, and will use a default location for the necessary MPI libraries.
- If -single is not used, then double precision executables will be created at the make [] command.
- Once installation is complete, a makefile will automatically be created for the machine platform on which the code is installed.
- Go to the build directory.
- By typing "make" you will see all the make options available.



- Several of the most common make options are:
 - make cfl3d seq make the sequential (single processor) version of the code
 - make cfl3d_mpi make the MPI (multiprocessor) version of the code
 - make splitter make the block splitter executable
 - make cfl3d_tools make some of the cfl3d utilities
- Within the build directory, type the make option for the executable you want.
- To execute the sequential code type:
 - ./cfl3d_seq < cfl3d.inp
- To execute the MPI code type:

mpirun –np <noprocessors> ./cfl3d_mpi < cfl3d.inp where <noprocessors> is typically one greater than the number of blocks*

^{*} The MPI process requires an extra administrative processor beyond those that perform the computation. (e.g. For a 12 block grid, all with equal numbers of grid points, to be run on 3 processors, noprocessors = 4)

Equations and Dimensions



Reference parameters

- The governing equations are the Euler or Navier-Stokes equations combined with a turbulence model for RANS computation
- The governing equations are non-dimensionalized based on the following parameters:

 $\widetilde{L}_{\scriptscriptstyle R}$ — Reference length used by the code (dimensional)

 $\widetilde{
ho}_{\!\scriptscriptstyle \infty}$ — Free-stream density, mass/unit length cubed

 $\widetilde{a}_{\cdot \cdot}$ — Free-stream speed of sound, length/time

 $\widetilde{\mu}_{\!\scriptscriptstyle \infty}$ — Free-stream molecular viscosity, mass/length-time

Equations and Dimensions



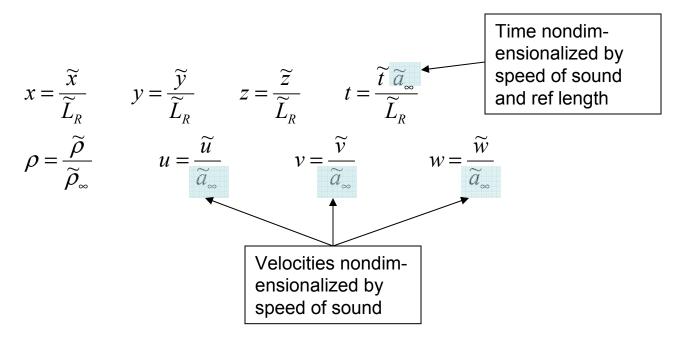
- Since there is no standard system of units for CFD models the non-dimensionalization in CFL3D removes the necessity of converting grids into units compatible with the code. The way in which the non-dimensionalization is accomplished will be presented later in this document.
- Note that the term free-stream is used in the non-dimensionalization. CFL3D was developed primarily as an external flow solver. It has the capability to perform computations for internal flows as well. Therefore a more general term reference state should probably be used, but the term free-stream is used throughout the documentation.

Equations and Dimensions



Non-dimensional variables

In CFL3D the non-dimensionalizations are performed as follows:



Non-dimensionalizing by speed of sound makes transonic the natural flow regime for CFL3D, although low speed and hypersonic flows can be computed, with modified input, as well.



Overview

- There are five steps in problem formulation and setup for steady and unsteady computation:
 - Condition definition
 - Grid generation
 - Block splitting (if necessary)
 - Blocking and boundary conditions
 - Input development
- Parameters that define a condition are:
 - Mach number
 - Reynolds number
 - Ambient temperature
 - Grid orientation (angle of attack, side slip, etc...)

Input for these parameters will be discussed later. For the moment several of these parameters are required for the proper construction of the grid...



Grid generation

Considerations that are important for generation of a grid:

- Reynolds number sets permissible Δy^{\dagger} at the surface.
 - For most turbulent computations typically want a $y^+ \sim 1$ for first grid off the surface
 - For turbulent computations with wall function, typically want a $y^+ \sim 50\text{-}100$ for first grid off the surface
 - Setting \(\Delta y^+\) requires an estimate of the wall shear stress prior to computing

Note that:
$$y^+ = \frac{y}{v} \sqrt{\frac{\tau_w}{\rho}}$$

where y is normal distance to surface, τ_w is wall shear stress, ρ is density and v is kinematic viscosity.



Grid generation

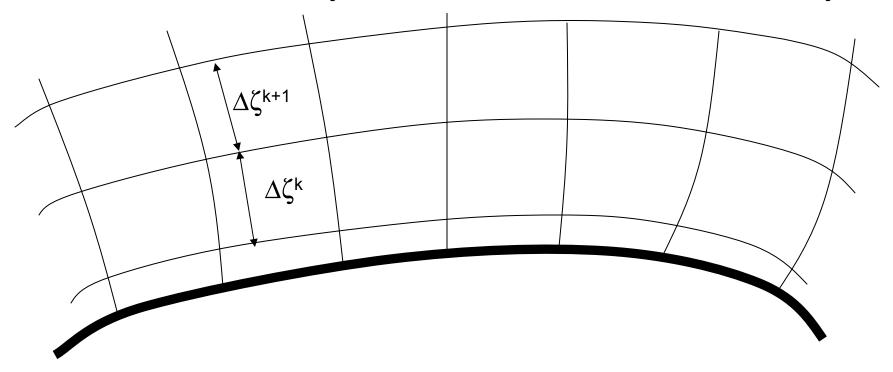
 After the first converged successful run with a coarse grid, y⁺ of the first grid can be checked. This value is found at the end of the cfl3d.out file. (See Y+ MAX, Y+ MIN and Y+AVG below)

```
YPLUS STATISTICS (endpts not included) - BLOCK 1 (GRID 1)
```

```
K=1 SURFACE:
 Y+ MAX JLOC ILOC
                     Y+ MIN JLOC ILOC
0.535E+00 151
                     0.261E-01 217
                1
 DN MAX JLOC ILOC
                      DN MIN JLOC ILOC
0.152E-05 228
                     0.149E-05 219
 Y+ AVG Y+ STD DEV
                      NY+>5 NPTS
0.264E+00 0.373E+00
                         0
                               199
YPLUS STATISTICS (endpts not included) - ALL GLOBAL BLOCKS
 Y+ MAX ILOC JLOC KLOC BLOCK GRID
0.535E+00 1
               151
                     1
 Y+ MIN ILOC JLOC KLOC BLOCK GRID
               217
0.261E-01
           1
     etc...
```

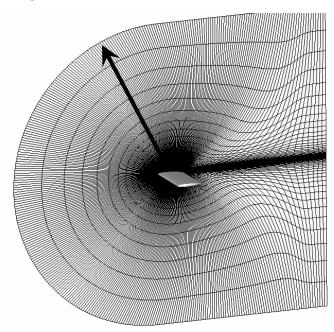


- Grid stretching away from a surface.
 - Rule of thumb: $\Delta \zeta^{k+1}$ should be no more than 1.2 to 1.5 times $\Delta \zeta^k$





- Outer extent of grid
 - Rule of thumb: The outer boundary of the grid should be at least 15 body lengths away (3D) and at least 20 body lengths away (2D). This is not a hard and fast rule and there are some notable exceptions. Note that the grid below would not be considered a fully acceptable grid.





- Grid quality
 - Grid metric smoothness. CFL3D assesses the size of local variations in grid metrics. Warnings are printed to the cfl3d.out file. Any messages of the following form indicate a problem with the grid:

```
FATAL si grid normal direction change near j,k,i,i+1= 23 5 164 165 ... suspect bad grid

FATAL sj grid normal direction change near j,k,i,i+1= 23 5 164 165 ... suspect bad grid

Etc... Or

WARNING: Dramatic si grid norm direction change (>120deg)

WARNING: Dramatic sj grid norm direction change (>120deg)

Etc...
```



- Grid quality (...continued)
 - Negative grid volumes. CFL3D checks whether there are negative volumes in the grid. Under normal operating procedures the code will exit with an error message in the cfl3d.error file.*
- Grid clustering to resolve flow gradients
 - Resolving a wake. Although angle of attack is specified in the input, it does result in the possibility of flow separation and wing stall and resulting wake. The wake may need grid clustering.
 - Resolving a shock or curvature effect. Mach number effects such as a shock or surface curvature may result in gradients that require resolving.
 - These steps must be performed prior to running CFL3D.

^{*} There is a keyword option that allows computing to continue with negative volumes. This option will be discussed later in the course under "Keyword Input".



Grid generation

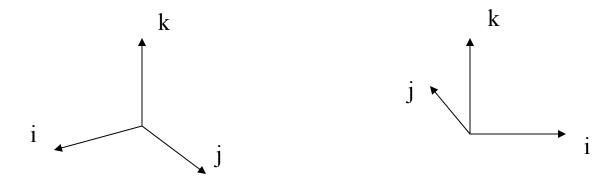
- Grid file format
 - The grid file format must be unformatted
 - Two grid data formats are possible, plot3d and cfl3d. These formats are presented in the CFL3D version 5.0 manual.
 - If CFL3D is compiled in double precision, the grid file must be written as double precision real
 - Example of multi-platform issue: If a Linux compiler is used to compile CFL3D to read an SGI unformatted grid file, the grid file must be generated with the same compile options

Example: Suppose the code 'hygrid' is used to generate the unformatted grid file. On a Linux based PC platform using the Portland Group compiler, the compile option –byteswapio swaps bytes from big-endian to little-endian for input compatibility with a Sun or SGI system. This compiler option will allow CFL3D to read the grid file created either on the PC cluster or on an SGI machine.



Grid generation

CFL3D requires that the right-hand rule be observed in both the x,y,z orientation and the i,j,k index directions. Also, i,j and k do not have to be in the x,y and z directions. Any permutation is valid as long as the right-hand rule is upheld. Caveat: When using turbulence models there are direction preferences as will be discussed.





Multigridable dimensions

To use multigrid, grid dimensions including all b.c. segments must be multigridable

Table 7-2. Grid sizes multigridable to three additional level.													
Grid:	Coarser Levels:				Grid:	Coarser Levels:			Grid:	Coa	Coarser Levels:		
9	5	3	2		345	173	87	44	673	337	169	85	
17	9	5	3		353	177	89	45	681	341	171	86	
25	13	7	4		361	181	91	46	689	345	173	87	
33	17	9	5		369	185	93	47	697	349	175	88	
41	21	11	6		377	189	95	48	705	353	177	89	
49	25	13	7		385	193	97	49	713	357	179	90	
57	29	15	8		393	197	99	50	721	361	181	91	
65	33	17	9		401	201	101	51	729	365	183	92	
73	37	19	10		409	205	103	52	737	369	185	93	
81	41	21	11		417	209	105	53	745	373	187	94	
89	45	23	12		425	213	107	54	753	377	189	95	
97	49	25	13		433	217	109	55	761	381	191	96	
105	53	27	14		441	221	111	56	769	385	193	97	
113	57	29	15		449	225	113	57	777	389	195	98	
121	61	31	16		457	229	115	58	785	393	197	99	
129	65	33	17		465	233	117	59	793	397	199	100	
137	69	35	18		473	237	119	60	801	401	201	101	
145	73	37	19		481	241	121	61	809	405	203	102	
153	77	39	20		489	245	123	62	817	409	205	103	
161	81	41	21		497	249	125	63	825	413	207	104	

From CFL3D User's Manual, 7.1.2, pg 129



Multigrid dimensions

				_							
169	85	43	22	505	253	127	64	833	417	209	105
177	89	45	23	513	257	129	65	841	421	211	106
185	93	47	24	521	261	131	66	849	425	213	107
193	97	49	25	529	265	133	67	857	429	215	108
201	101	51	26	537	269	135	68	865	433	217	109
209	105	53	27	545	273	137	69	873	437	219	110
217	109	55	28	553	277	139	70	881	441	221	111
225	113	57	29	561	281	141	71	889	445	223	112
233	117	59	30	569	285	143	72	897	449	225	113
241	121	61	31	577	289	145	73	905	453	227	114
249	125	63	32	585	293	147	74	913	457	229	115
257	129	65	33	593	297	149	75	921	461	231	116
265	133	67	34	601	301	151	76	929	465	233	117
273	137	69	35	609	305	153	77	937	469	235	118
281	141	71	36	617	309	155	78	945	473	237	119
289	145	73	37	625	313	157	79	953	477	239	120
297	149	75	38	633	317	159	80	961	481	241	121
305	153	77	39	641	321	161	81	969	485	243	122
313	157	79	40	649	325	163	82	977	489	245	123
321	161	81	41	657	329	165	83	985	493	247	124
329	165	83	42	665	333	167	84	993	497	249	125
337	169	85	43								

From CFL3D User's Manual, 7.1.2, pg 129



Blocking and boundary conditions

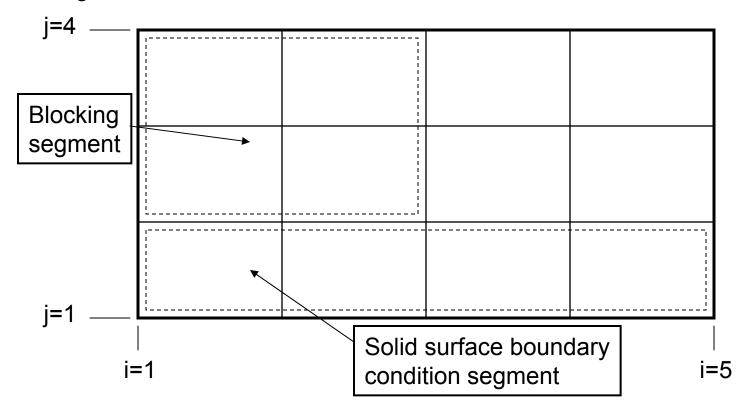
Blocking and boundary conditions are specified at the following boundaries:

where idim, jdim and kdim are the block dimensions in the ijk-directions. Blocking and boundary condition data can be composed of multiple segments but the combined segments must span each of the six block faces. Note that to perform multigrid computations, the boundary and blocking segments must be multigridable integers.



Blocking and boundary conditions

Example of possible blocking or boundary condition segments on the k0 face. Suppose that part of the k0 face below represents the surface of a wing.

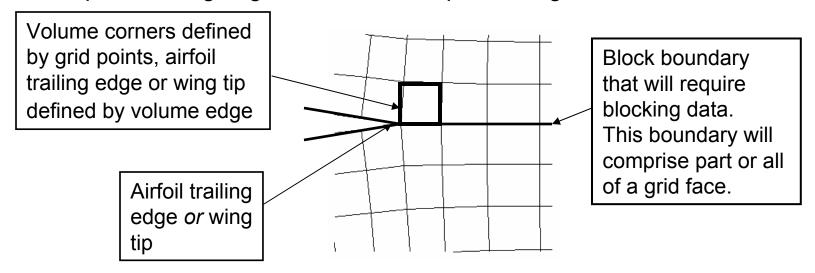




Blocking and boundary conditions

Volume *edges* define geometric extremities. The volume edges will also be the start and end points of blocking pairs. All blocking and boundary conditions will be on external surfaces of grid blocks.

Example: Trailing edge of an airfoil or tip of a wing.

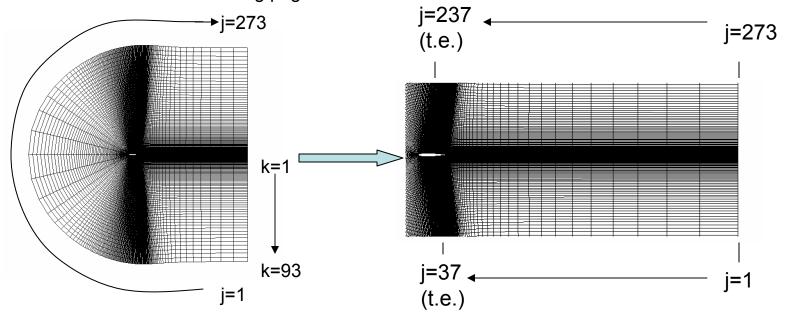




Blocking and boundary conditions

Blocking defines the start and ending indices of 1-1 interfaces between one or more corresponding grid blocks.

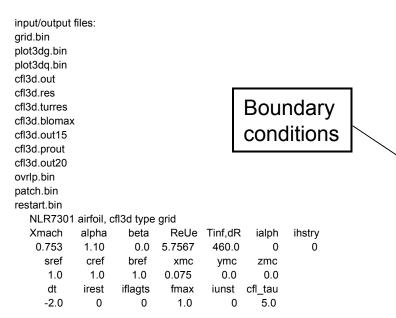
Consider the example of a 2D airfoil using a single block C-grid with dimension 2x273x93. CFL3D is a finite volume code and therefore requires 2 grid points in the span-wise direction (always i-dir for a 2D grid). *Note that the arrows in the right hand figure below denotes the end of a blocking segment.* The meaning of this statement will be made clear in the following pages.





Blocking and boundary conditions

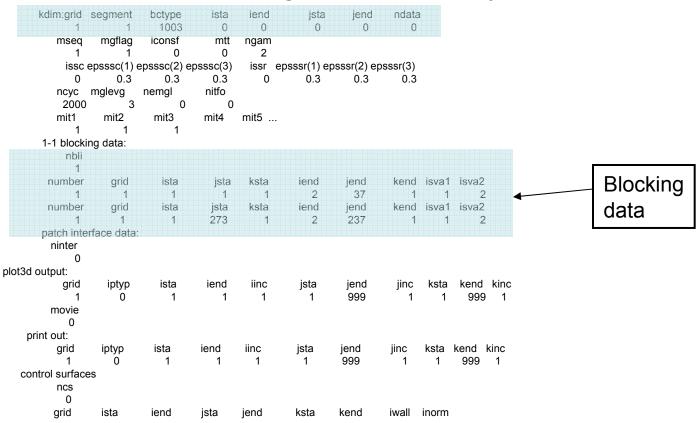
The following is the steady input file for the single block C-grid 2D airfoil. Highlighted sections are the blocking and boundary condition input:



ngrid	nplot3d		nwrest			ntstep	ita	
1	1	1	1000		1	1	-2	
ncg		iadvance	iforce	ivisc(i)	ivisc(j)	ivisc(k)		
2	0	0	1	0) () 5		
idim	jdim	kdim						
2	273	93						
ilamlo		jlamlo	-		klamhi			
0	0	0	0	0	0			
inewg	igridc		js	ks	ie	je	ke	
0	0	0	0	0	0	0	0	
idiag(i)	idiag(j)	idiag(k)	iflim(i)	iflim(j)	iflim(k)			
1	1	1	4	4	4			
ifds(i)	ifds(j)	ifds(k)		rkap0(j)				
1	1	1		0.3333				
grid	nbci0	nbcidim	nbcj0	nbcjdim		nbckdim	iovrlp	
1	1	1	1	1	3	1	0	
i0: grid	segment			jend	ksta		ndata	
1	1	1002		0	0	0	0	
idim:grid	segment	bctype			ksta			
1	1	1002		0	0	0	0	
j0: grid	segment	bctype		iend	ksta	kend	ndata	
1	1	1003	0	0	0	0	0	
jdim:grid	segment	bctype	ista	iend	ksta	kend	ndata	
1	1	1003	0	0	0	0	0	
k0: grid	segment	bctype	ista	iend	jsta	jend	ndata	
1 1	0	0	0	0	1	37	0	
1 2	2004	0	0	0	37	237	2	
tw/tinf	cq							
0.	0.							
0.	3				237	273		



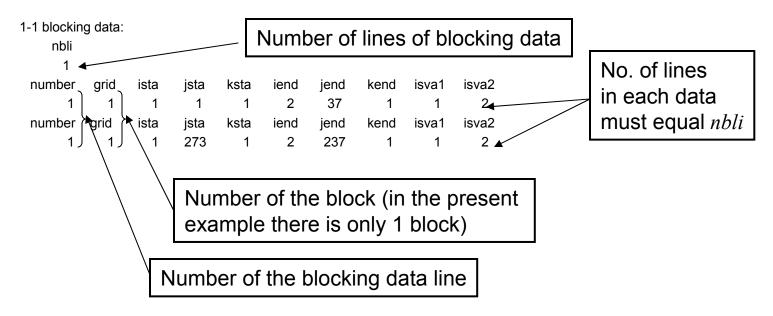
Blocking and boundary conditions





Blocking and boundary conditions

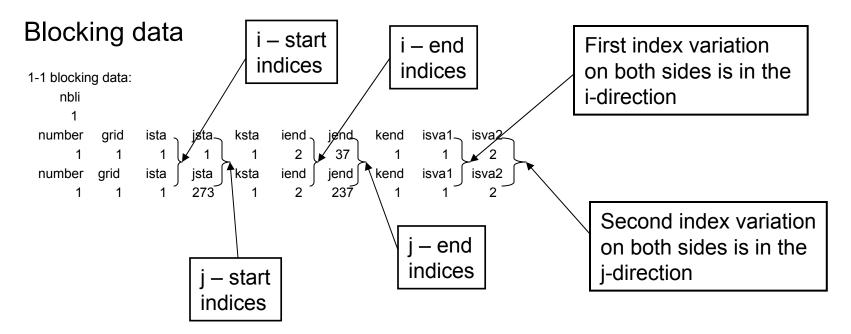
For this example, format of the blocking data in the input file:



Note: The text cards must be present, but the text within those lines is arbitrary, and is for user information only. All lines with data are in free field format throughout the input file.



Blocking and boundary conditions

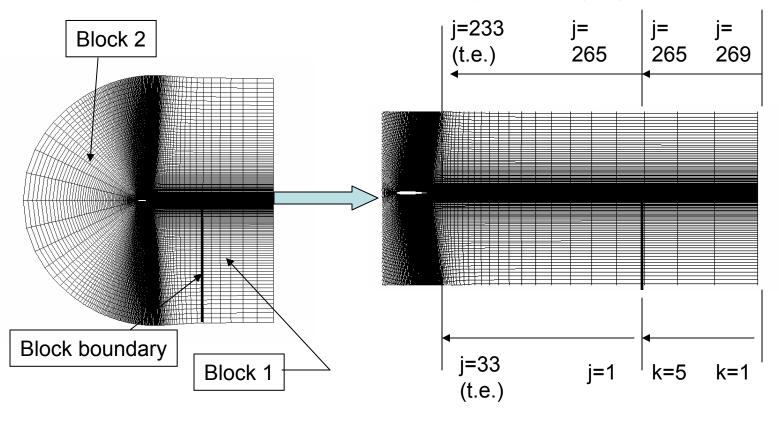


Because this is a volume grid, the blocking will always define a two-dimensional interface in index space



Blocking and boundary conditions

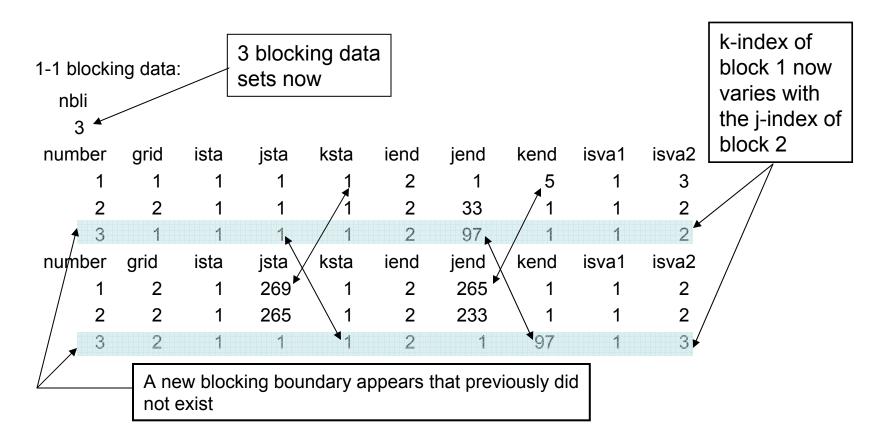
Consider a second example of a 2D airfoil using two blocks to compose a C-grid. Block 1 has dimensions 2x93x5. Block 2 has dimensions 2x269x93. Note again that the arrows in the right hand figure below denotes the end of a blocking segment. This fact is made clear by the following page.





Blocking and boundary conditions

Blocking data

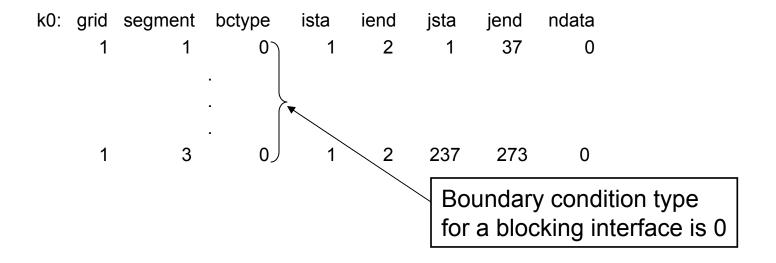




Blocking and boundary conditions

Blocking faces require corresponding boundary condition data

In the first example above, the blocking interface is at the k=1 boundary. Therefore, the boundary condition data for that blocking interface is in the 'k0' boundary data.

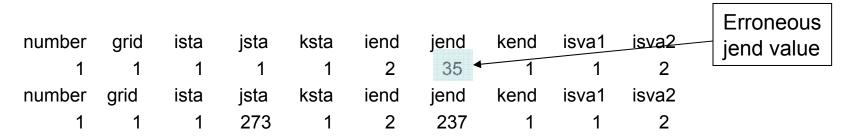




Blocking and boundary conditions

CFL3D will stop if the number of grid points across a blocking interfaces does not match.

Suppose the following blocking data had been specified for example 1 above:



Execution will terminate with the following error message at the end of the file 'precfl3d.out':

.

the limits of ind2 are not the same for both sides for 1:1 plane 1



Blocking and boundary conditions

CFL3D also checks the input connection data by computing the geometric mismatch between both sides of the interface. A true 1-1 interface will have zero (machine zero) mismatch. Any mismatches larger than ε (where ε is the larger of 10⁻⁹ or 10x(machine zero)) will cause a warning message.

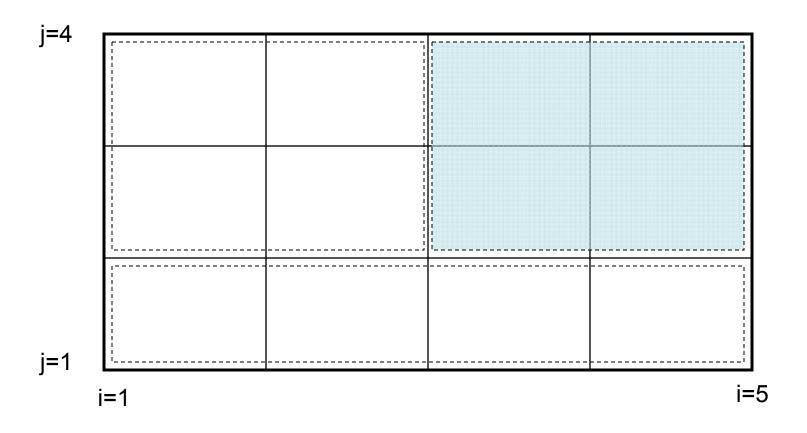
Example of the output in 'cfl3d.out':

```
j= 1 1-1 blocking type 0 i= 1, 31 k=137, 69 connects to j = 1 of block 2 blocking check....geometric mismatch = 0.2166272E-03
```



Blocking and boundary conditions

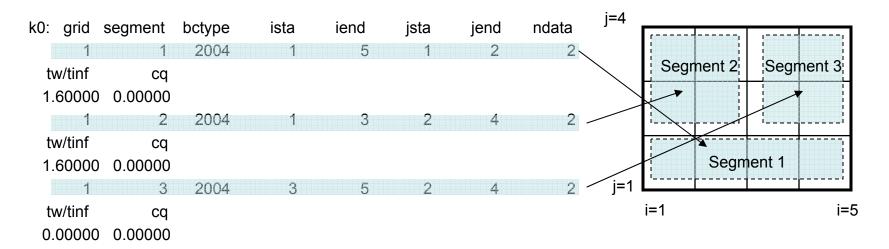
Example of possible boundary condition segments on the k0 face. Suppose that the k0 face below represents the surface of a wing.





Blocking and boundary conditions

At the unshaded cells, it is desired to apply a heated wall boundary condition, while at the shaded cells it is desired to apply an adiabatic wall boundary condition. One way to accomplish this objective is to divide the boundary into the segments shown. The CFL3D input file would have input that looks like this:



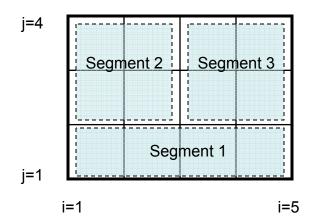
Note that for segment 1, for instance, the grid points i = 1 to 5, j = 1 to 2 define the boundary of the cells at which the condition type is to be applied.



Blocking and boundary conditions

Setting ista = iend = 0 and/or jsta = jend = 0 is a shorthand way of specifying the entire range. In other words, an alternate boundary condition input with identical outcome is:

k0:	grid	segment	bctype	ista	iend	jsta	jend	ndata
	1	1	2004	0	0	1	2	2
tv	v/tinf	cq						
1.6	60000	0.00000						
	1	2	2004	1	3	2	4	2
tv	v/tinf	cq						
1.6	60000	0.00000						
	1	3	2004	3	5	2	4	2
tv	v/tinf	cq						
0.0	00000	0.00000						





Blocking and boundary conditions

The following 1000 series boundary conditions are available:

bctype	boundary condition
1000	free stream
1001	general symmetry plane
1002	extrapolation
1003	inflow/outflow
1005	inviscid surface
1006	inviscid surface (using normal momentum)
1008	tunnel inflow
1011	singular axis – half-plane symmetry
1012	singular axis – full plane
1013	singular axis – partial plane

Refer to the Version 5.0 Manual and Version 6.0 web page for more information on these boundary conditions



Blocking and boundary conditions

The following 2000 series boundary conditions are available:

bctype	boundary condition
2002	specified pressure ratio
2003	inflow with specified total conditions
2004	no-slip wall
2005	periodic in space
2006	set pressure to satisfy the radial equilibrium equation
2007	set all primitive variables

Refer to the Version 5.0 Manual and Version 6.0 web page for more information on these boundary conditions



Blocking and boundary conditions

The following 2000 series boundary conditions are available:

bctype	boundary condition
2008	user specifies density and velocity components, pressure extrapolated from interior
2009	sets total pressure and total temperature. Inflow pressure
	extrapolated from interior
2014	user specifies transpiration through the boundary
2018	user specifies temperature and momentum components, pressure extrapolated from interior
2028	user specifies frequency and maximum momentum components, density and pressure extrapolated
2102	pressure ratio specified as a sinusoidal function of time

Refer to the Version 5.0 Manual and Version 6.0 web page for more information on these boundary conditions



Blocking and boundary conditions

Boundary condition 1000 - Free stream. Extrapolation points just outside the boundary are set to initial free stream values, which are:

$$\rho_{initial} = 1.0$$

$$u_{initial} = M_{\infty} \cos \alpha \cos \beta$$

$$v_{initial} = -M_{\infty} \sin \beta$$

$$w_{initial} = M_{\infty} \sin \alpha \cos \beta$$

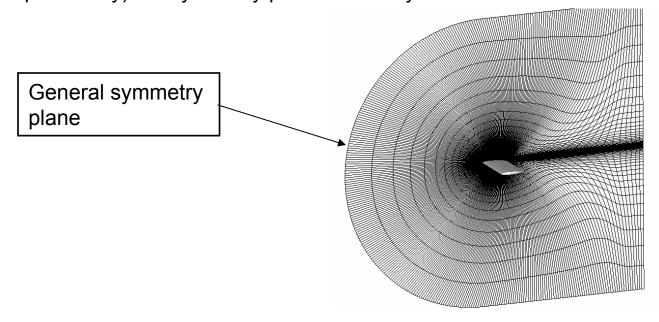
$$p_{initial} = \rho_{initial} a_{initial}^{2} / \gamma$$

where ρ is density, u,v,w are the x,y,z components of velocity, p is pressure, a is speed of sound and γ is ratio of specific heats. M is Mach number, α is the angle of attack and β is the side slip angle.



Blocking and boundary conditions

Boundary condition 1001 - General symmetry plane. Suppose we wish to simulate a 3D wing using the half wing shown. If only one type of maneuver is performed (i.e. with aircraft maneuver symmetry in the x-y plane, x-z plane or y-z plane only) the symmetry plane boundary condition can be used.





Blocking and boundary conditions

Boundary condition 1002 - Extrapolation. Ghost points outside the flow field domain are extrapolated from the interior.

Boundary condition 1003 - Inflow/Outflow. This condition uses Riemann invariants to calculate inflow and outflow at the boundary cell face. It effectively sets total pressure.

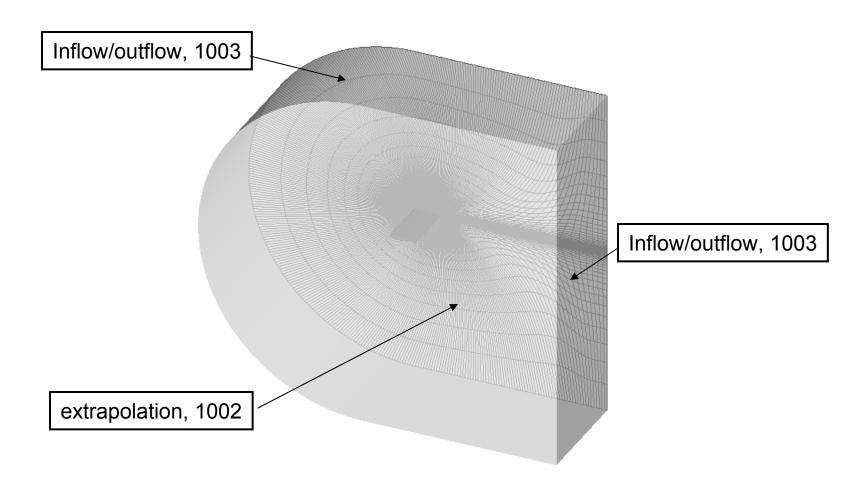
Boundary condition 1005 - Inviscid surface. Velocity components normal to the surface are set to zero. Density and pressure gradients are set to zero.

Boundary condition 1006 - Inviscid surface. Similar to b.c. 1005 except that the normal momentum equation is used to obtain wall pressure. Generally results in a smoother solution near an inviscid surface.

Boundary condition 2004 - No slip wall. Viscous boundary conditions are set at surface cell face, i.e. flow velocity equals the surface velocity.



Example of typical "outer" boundary conditions





Blocking and boundary conditions

Boundary condition 1005: Inviscid surface

•

i0:	grid	segment	bctype	ista	iend	jsta	jend	ndata
	1	1	1005	1	5	1	2	0
	1	2	0	1	3	2	4	0
idim	n:grid	segment	bctvpe	ista	iend	ista	iend	ndata

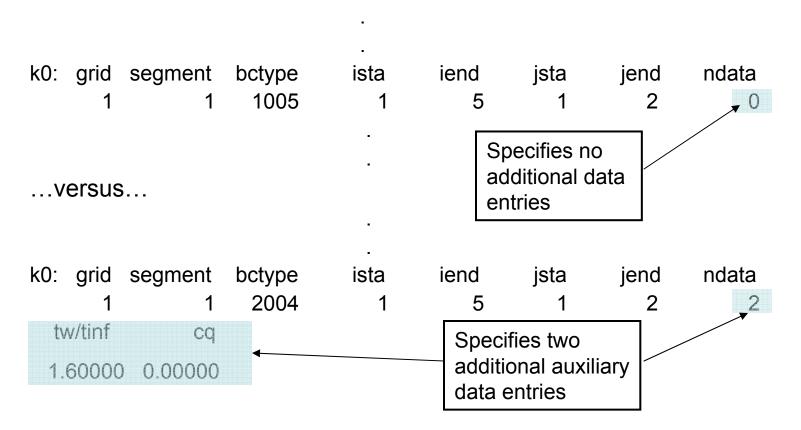
.

.



Blocking and boundary conditions

Note that the b.c. 1005 has no auxiliary data, while the b.c. 2004 has two additional lines





Blocking and boundary conditions

- Series 1000 boundary conditions require no auxiliary data
- Number of auxiliary data entries for series 2000 boundary conditions are shown below

No. of auxiliary
data
1
5
2
5
4
5*
4*
4*
3
7
4*
4*
4

^{*} Means turbulence data can also be specified, adding either 1 or 2 additional aux. data inputs

See the CFL3D version 5.0 manual and CFL3D Version 6 web page for discussion of these boundary conditions



Blocking and boundary conditions

Example of a boundary condition with 5 auxiliary data entries: 2003 - "Engine inflow", inflow with specified total conditions:

•

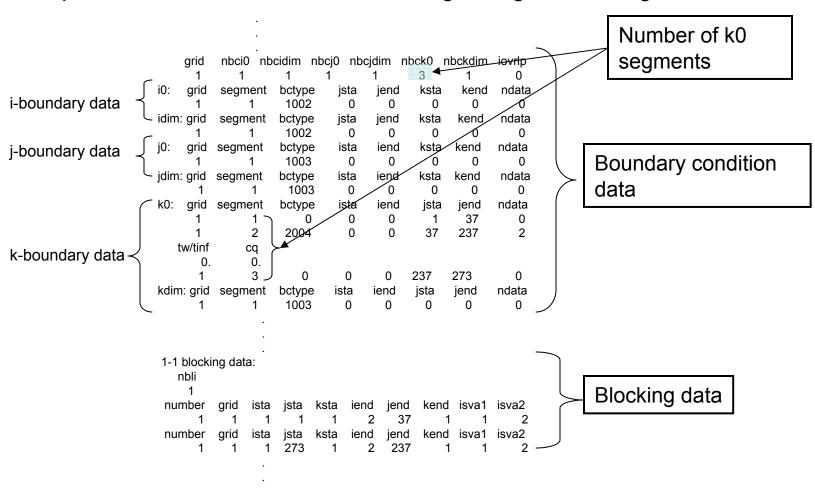
k0:	grid	segment	bctype	ista	iend	jsta	jend	ndata
	1	1	2003	1	5	1	2	5
M	lach	Pt/Pinf	Tt/Tinf	Alphae	Betae			
	0.30	4.000	1.1755	0.0	0.0			

.



Blocking and boundary conditions

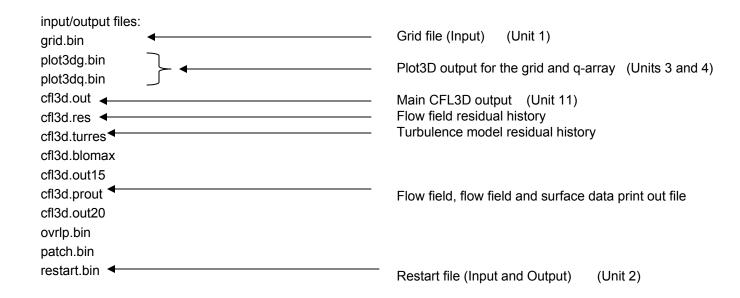
Input data so far for the 2D airfoil using a single block C-grid





Input/output file specifications

Some of the key input, output files:





Input/output file specifications

- These names can be changed by the user.
- Input/output redirects are permitted. (e.g. ../../grid.bin or ./cflout/cfl3d.out)
- Additional files are printed out not contained in this list. (e.g. precfl3d.out, precfl3d.error, cfl3d.error, cfl3d.subit_res and cfl3d.subit_turres)
- The restart file name that is read at the start of the computation is the same name used for output at the end. Scripting that saves restart files to another name will be required if the user wishes to save the input restart.



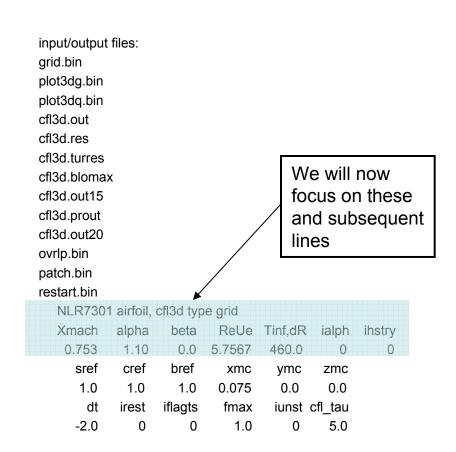
Navigating diagnostic output

Diagnostic output:

- Initial input syntax and completeness are checked in the preprocessor 'precfl3d'. This is an initial step automatically performed by CFL3D. Output from this check will be in the files 'precfl3d.error' and 'precfl3d.out'. Input errors will cause the output in 'precfl3d.out' to stop at the line at which the error occurred. Often informative diagnostics will be output there.
- When the checker 'precfl3d' has determined that the input is properly configured, the top of 'cfl3d.out' will show the input values it has read.
- Other checks (e.g. grid dimension, blocking, incompatibility of a restart file) are performed in 'cfl3d'. Error output including the suspected cause of the termination will be found in 'cfl3d.error'. Sometimes additional insight into the cause of the error can be found by checking the main output in 'cfl3d.out' although frequently there is little additional diagnostic output in 'cfl3d.out' if the code terminates.

NASA

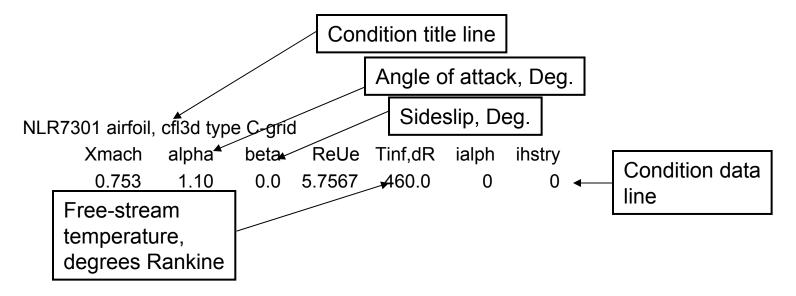
Title line and condition data



	ngrid	nplot3d	nprint 1	nwrest 1000	ichk 0	i2d	ntstep	ita -2
	nca	iom	iadvance	iforce	ivisc(i)	ivisc(j)	ivisc(k)	-2
	ncg 2	0	nauvance 0	1	0 (1)		1VISC(K)	
	idim	idim	kdim	'	U	U	5	
	2	273	93					
	ilamlo			jlamhi	klamlo	klamhi		
	0	0	0	0	0	0		
	inewa	igridc	is	js	ks	ie	je	ke
	0	.9.140	0	0	0	0	0	0
i	diag(i)	idiag(j)	•	iflim(i)	•	iflim(k)	·	ŭ
•	1	1	1	4	4	4		
	ifds(i)	ifds(j)	ifds(k)	rkap0(i)	rkap0(j)	rkap0(k)		
	1	1	1	0.3333		0.3333		
	arid	nbci0	nbcidim				nbckdim	iovrlp
	1	1	1	1	1	3	1	Ö
i0:	arid	segment	bctype	ista	iend	ksta	kend	ndata
	1	1	1002	0	0	0	0	0
idi	m:grid	segment	bctype	jsta	jend	ksta	kend	ndata
	1	1	1002	0	0	0	0	0
j0:	grid	segment	bctype	ista	iend	ksta	kend	ndata
-	1	1	1003	0	0	0	0	0
jdi	m:grid	segment	bctype	ista	iend	ksta	kend	ndata
	1	1	1003	0	0	0	0	0
k0	: grid	segment	bctype	ista	iend	jsta	jend	ndata
1	1	0	0	0	0	1	37	0
1	2	2004	0	0	0	37	237	2
	tw/tinf	cq						
	0.	0.						
	1	3	0	0	0	237	273	0



Title line and condition data



ialph – indicator to determine whether angle of attack is measured in the x-z plane or the x-y plane ihstry – determines which variables are to be tracked for convergence history. Default is $C_{\rm l}$, $C_{\rm d}$, $C_{\rm y}$ (or $C_{\rm z}$), $C_{\rm m.}$

Input of ReUe (Reynolds number) requires some additional explanation....



Reference length

Calculation of ReUe

Recall the nondimensionalizations:

$$x = \frac{\widetilde{x}}{\widetilde{L}_{R}} \qquad y = \frac{\widetilde{y}}{\widetilde{L}_{R}} \qquad z = \frac{\widetilde{z}}{\widetilde{L}_{R}} \qquad t = \frac{\widetilde{t} \ \widetilde{a}_{\infty}}{\widetilde{L}_{R}}$$

$$\rho = \frac{\widetilde{\rho}}{\widetilde{\rho}_{\infty}} \qquad u = \frac{\widetilde{u}}{\widetilde{a}_{\infty}} \qquad v = \frac{\widetilde{v}}{\widetilde{a}_{\infty}} \qquad w = \frac{\widetilde{w}}{\widetilde{a}_{\infty}}$$

Reynolds number based on reference length:

$$\operatorname{Re}_{\widetilde{L}_{R}} = \frac{\widetilde{\rho}_{\infty} |\widetilde{V}_{\infty}| \widetilde{L}_{R}}{\widetilde{\mu}_{\infty}}$$



Calculation of ReUe

Calculation of ReUe

$$ReUe = Re_{\widetilde{L}_R} \times 10^{-6} = \frac{\widetilde{\rho}_{\infty} |\widetilde{V}_{\infty}| \widetilde{L}_R}{\widetilde{\mu}_{\infty}} \times 10^{-6} = \frac{\widetilde{\rho}_{\infty} M_{\infty} \sqrt{\gamma R T_{\infty}} \widetilde{L}_R}{\widetilde{\mu}_{\infty}} \times 10^{-6}$$

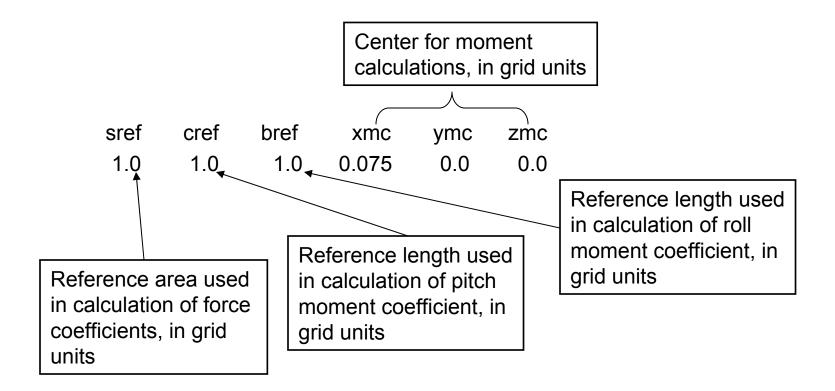
Example: Suppose we have a grid that is in inches, and we wish to retain that length scale so that the grid remains compatible with a finite element model of the wing structure that is also in inches. Suppose the Reynolds number is 1 million based on chord length of 20 inches.

Set
$$\widetilde{L}_{R}=1$$
 inch , then $\operatorname{Re}_{\widetilde{L}_{R}}=\operatorname{Re}_{c}(\widetilde{L}_{R}/c)=50{,}000,$ $\operatorname{Re}Ue=0.05$

ReUe is the Reynolds number per unit grid length in millions

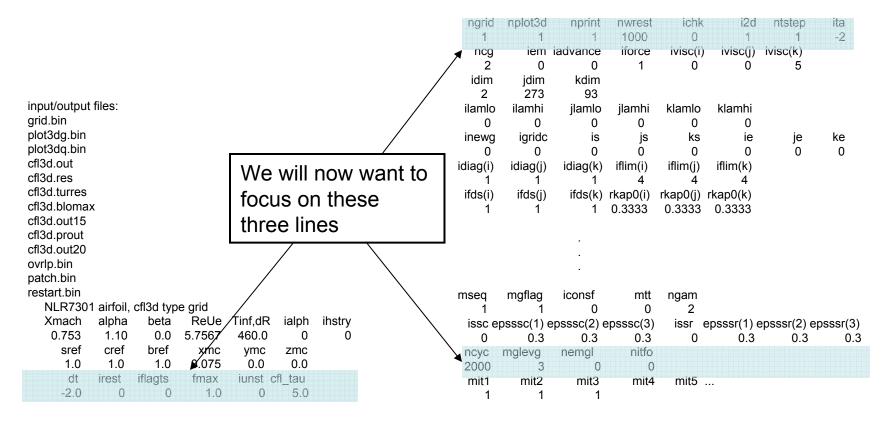
Reference data input





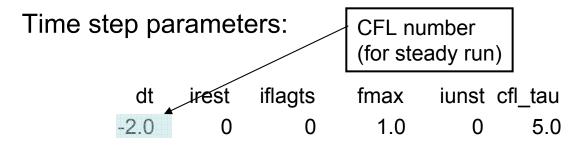


Steady solution cycling input

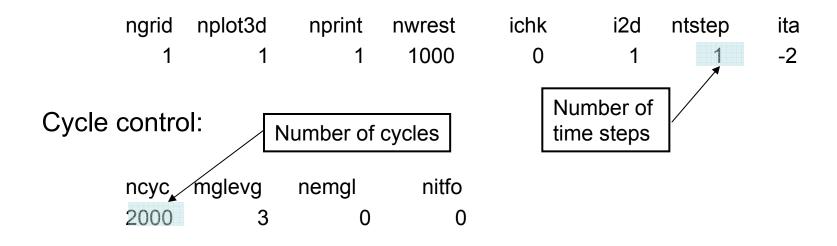




Steady solution cycling input



Number of time step advances, and time accuracy:





Steady solution cycling input

dt		iflagts	fmax	iunst	cfl_tau		
-2.0	0	0	1.0	0	5.0		
ngrid	nplot3d	nprint	nwrest	ichk	i2d	ntstep	ita -2
ı	I	. 1	1000	0	ı	ı	-2
ncyc	mglevg	nemgl	nitfo				
2000	3	0	0				

Note:

- when dt < 0, local time stepping is used, i.e. CFL = |dt|. This is used for converging a steady state solution. For steady state computations

$$\Delta \tau = CFL \cdot \Delta r$$

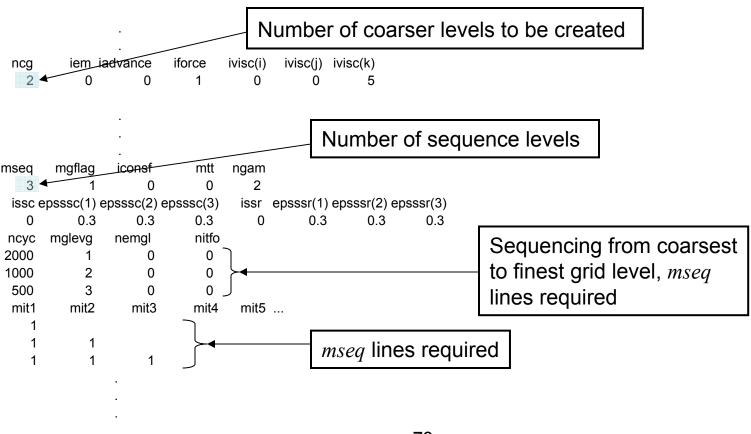
where Δr is a measure of local grid spacing and $\Delta \tau$ is the local pseudo time step size.

- cfl tau is not used when dt < 0. The value input for that parameter is a placeholder.
- *iunst* is set to 0 in the code when dt < 0.
- *ntstep* is set to 1 in the code when dt < 0.
- ncyc controls the number of steady solution cycles computed.
- Values of dt of -2.0 to -10.0 are typical. Lower values will be required for a stiffer problem.



Grid sequencing

Grid sequencing can and should be used to accelerate convergence to a steady state solution. The following input sequences through three grid levels.



NASA

Grid sequencing output

The following grid level information will be found in the cfl3d.out on the completion of the 3D single block C-grid airfoil computation:

Because ncg = 2, two coarser levels created

```
reading grid 1 of dimensions (I/J/K): 2 273 93
creating coarser block 2 of dimensions (I/J/K): 2 137 47
creating coarser block 3 of dimensions (I/J/K): 2 69 24

******** BEGINNING TIME ADVANCEMENT, iseq = 1 ******
steady-state computations

******** BEGINNING MULTIGRID CYCLE ******

iseq= 1
level top = 1
level bottom = 1
number of global grid levels = 1
lglobal= 1

Coarsest to mid level
```

```
****** BEGINNING SEQUENCING TO FINER LEVEL *****

interpolating solution on coarser block 3 to finer block 2 (grid 1)
jdim,kdim,idim (finer grid)= 137 47 2
jj2,kk2,ii2 (coarser grid)= 69 24 2
interpolating turb quantities from coarser to finer block

****** ENDING SEQUENCING TO FINER LEVEL *****

****** BEGINNING TIME ADVANCEMENT, iseq = 2 *****

steady-state computations

****** BEGINNING MULTIGRID CYCLE *****

iseq= 2
level top = 2
level bottom = 1
number of global grid levels = 2
lglobal= 2
.
```



Grid sequencing output

```
****** BEGINNING SEQUENCING TO FINER LEVEL ******

interpolating solution on coarser block 2 to finer block 1 (grid 1)

jdim,kdim,idim (finer grid)= 273 93 2

jj2,kk2,ii2 (coarser grid)= 137 47 2

interpolating turb quantities from coarser to finer block

****** ENDING SEQUENCING TO FINER LEVEL *****

****** BEGINNING TIME ADVANCEMENT, iseq = 3 *****

steady-state computations

******* BEGINNING MULTIGRID CYCLE *****

iseq= 3

level top = 3

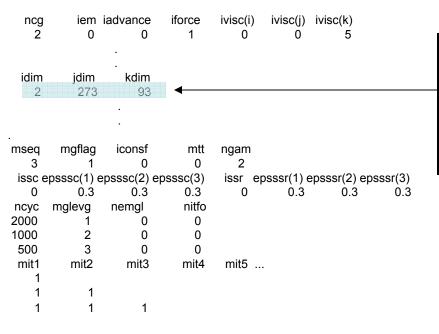
level bottom = 1
```

number of global grid levels = 3

Iglobal= 3



Grid sequencing



These dimensions support up to four multigrid levels. See version 5.0 manual for a table of multigridable dimensions. Note that *idim* is not multigridded for a 2D grid.

Note:

- The number of grid levels that will have been created are the coarser levels (ncg) plus the finest level. Therefore, mseq must be equal to or less than ncg + 1. Setting mseq higher than this will result in a termination and an error message in precfl3d.out.
- The permissible value of ncg will depend on the dimensions of the grid. It is usually good to have three to four possible levels of multi-grid. For example, since four levels of multi-grid are possible with this grid, we could have set ncg = 3.



Grid sequencing

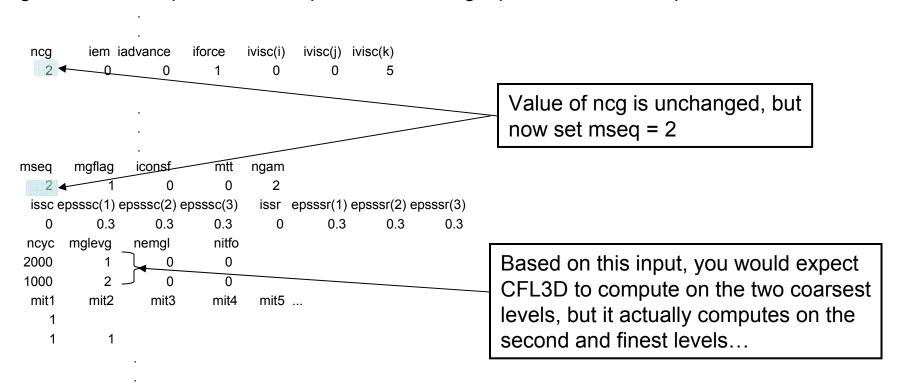
Note:

- Many more cycles will be done at the coarser levels. The computing required for a 3D grid will be a factor of 8 cheaper at each coarser level. For the present problem, the coarsest level would be 64 times cheaper than the finest level if a 3D grid had been used. Since it is a 2D grid it will be 16 times cheaper.
- It is usually good to completely converge the coarser levels before proceeding to the finer level. However, some problems will not compute well at a coarse level, but will compute at a finer level.
- Mglevg is always starting from the finest level ... as the following example will show...



Grid sequencing

Example: We wish to compute on only the two coarser levels with the grid used in the previous example. The following input has been set up:





Grid sequencing

...Here is what is actually output in cfl3d.out:

```
***** BEGINNING TIME ADVANCEMENT, iseq = 1 *****
steady-state computations
***** BEGINNING MULTIGRID CYCLE *****
iseq= 1
level top = 2
level bottom = 2
number of global grid levels = 1
Iglobal= 2
***** BEGINNING SEQUENCING TO FINER LEVEL *****
interpolating solution on coarser block 2 to finer block 1 (grid 1)
 jdim,kdim,idim (finer grid)= 273 93 2
 jj2,kk2,ii2 (coarser grid)= 137 47 2
 interpolating turb quantities from coarser to finer block
***** ENDING SEQUENCING TO FINER LEVEL *****
***** BEGINNING TIME ADVANCEMENT, iseq = 2 *****
```

```
steady-state computations

***** BEGINNING MULTIGRID CYCLE *****

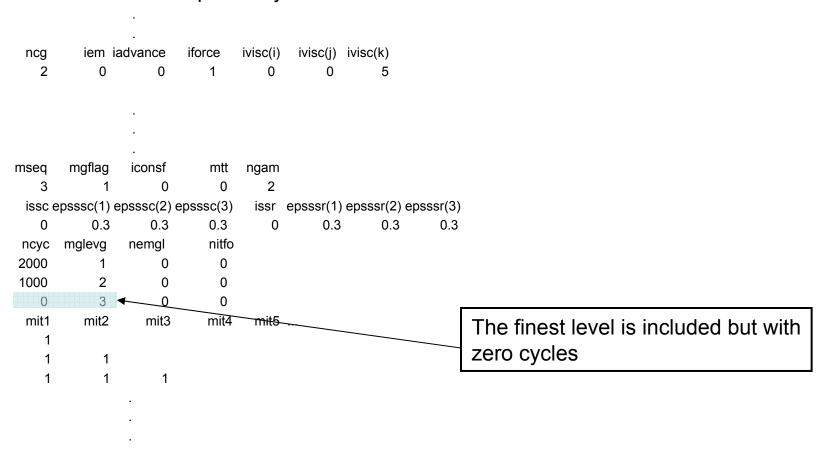
iseq= 2
level top = 3
level bottom = 2
number of global grid levels = 2
lglobal= 3
```

Computations performed on the middle and finest grids



Grid sequencing at coarsest levels only

Here is how to compute only on the two coarsest levels:



80



Grid sequencing at coarsest levels only

....and here is the output:

```
steady-state computations
***** BEGINNING TIME ADVANCEMENT, iseq = 1 *****
                                                                                 ***** BEGINNING MULTIGRID CYCLE *****
steady-state computations
                                                                                 iseq= 2
***** BEGINNING MULTIGRID CYCLE *****
                                                                                 level top = 2
                                                                                 level bottom = 1
iseq= 1
                                                                                 number of global grid levels = 2
level top = 1
                                                                                 Iglobal= 2
level bottom = 1
number of global grid levels = 1
Iglobal= 1
***** BEGINNING SEQUENCING TO FINER LEVEL *****
                                                                                  Computations performed on the
interpolating solution on coarser block 3 to finer block 2 (grid 1)
                                                                                  coarsest and middle levels
 jdim,kdim,idim (finer grid)= 137 47 2
 jj2,kk2,ii2 (coarser grid)= 69 24 2
 interpolating turb quantities from coarser to finer block
***** ENDING SEQUENCING TO FINER LEVEL *****
***** BEGINNING TIME ADVANCEMENT, iseq = 2 *****
```



Grid sequencing at coarsest levels only

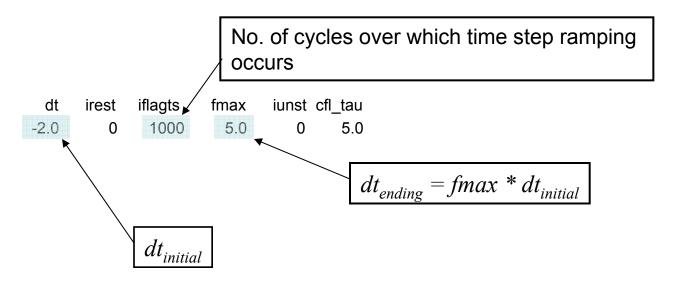
Why is it sometimes valuable to compute on the coarser levels only?

- Cost effectiveness of coarser levels
- Sometimes it is not possible to converge the finest level
- Many times you will want to compute unsteady solutions on coarser levels only, especially when debugging. Computing unsteady solutions on coarser levels only requires the steady starting point be on a coarser level.



Ramping up dt

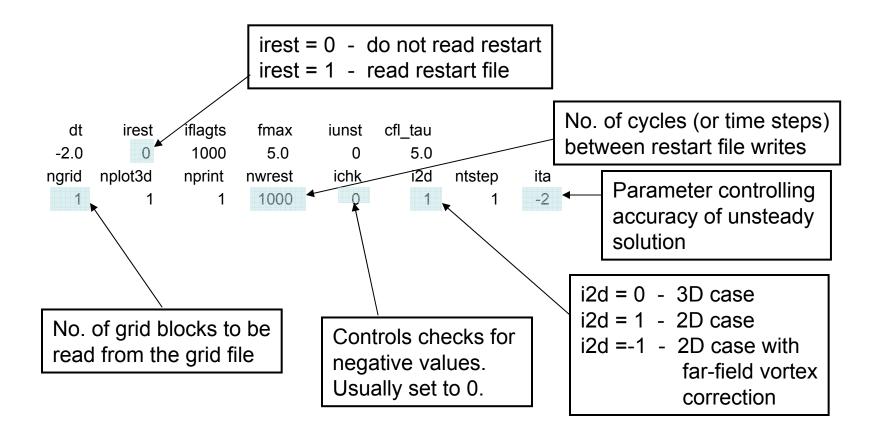
Sometimes it is useful for stiff problems to ramp up the time step size. Ramping up the time step size is accomplished with the following input:



In this example, the final CFL value of 10 is obtained after 1000 cycles. Note that this counter is reset with each restart. Therefore, $dt_{initial}$ will have to be reset to the dt_{ending} of the previous run.

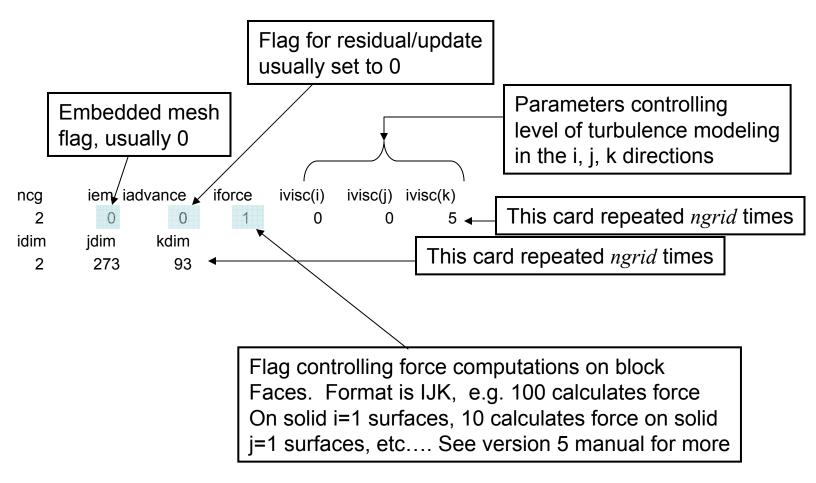
NASA

Additional input



NASA

Additional input



NASA

Turbulence model input

There are more than 13 turbulence models available, but the following are the most common turbulence models and the corresponding parameter input values:

0 - inviscid	
1 - laminar	
turbulent, Baldwin-Lomax with Degani- option (not recommended)	Schiff
5 - turbulent, Spalart-Allmaras model	
6 - turbulent, Wilcox k-ω	
7 - turbulent, k-ω SST (Menter's version)	
13 - nonlinear EASM k-ε model	
- nonlinear EASM k-ω model	

See the CFL3D Version 5.0 manual (Appendix H) and the CFL3D Version 6 web page (under `New Features') for descriptions of these and other models. See also under the 'Keywords' discussion in these notes for parameters that turn turbulence model features on.



Turbulence model

Several key notes on turbulence models:

- 1. If ivisc(m) < 0, a wall function is employed
- 2. Thin-layer viscous terms (laminar or turbulent) can be included in the i,j or k directions separately or combined. Cross-derivatives are not included. For the Baldwin-Lomax model, terms are allowed simultaneously in two directions only, either j-k or i-k.
- 3. Using the Baldwin-Lomax model with multi-zonal grids, wall distances are calculated only within a given zone.
- 4. It is preferable to let k be the primary viscous direction and i be secondary viscous direction.
- 5. The minimum distance function *smin* is computed from viscous walls only, not inviscid walls.

NASA

Turbulence model

- 6. Note that the field equation turbulence models may or may not transition to turbulent flow. Whether they transition will largely be determined by the free stream value of turbulence. Free stream turbulence level can be set in the key word input.
- 7. There are several places in which the turbulence level can be checked
 - There is an option allows the output of turbulence quantities in the plot3d file.
 - The file 'cfl3d.prout' contains the value of the turbulent viscosity. This is shown in the next slide.

See the CFL3D User's Manual, Version 5.0, Section 3.7 for more complete discussion



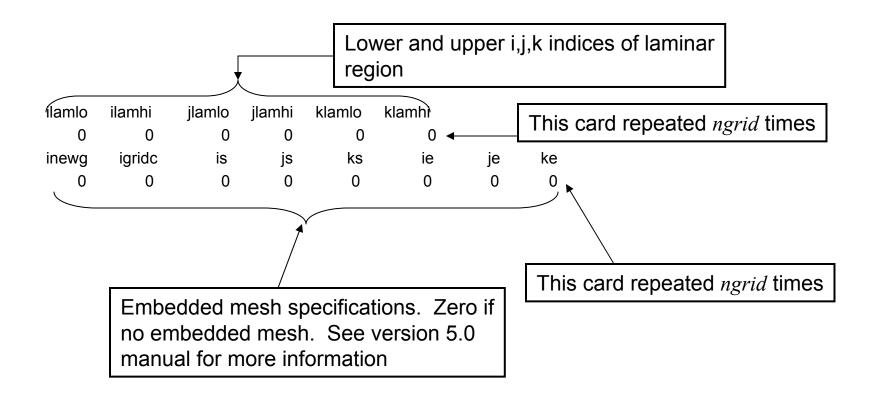
Turbulence model output



Data lines will be printed out for all flow field points specified by the user in the 'print out' portion of the input file.

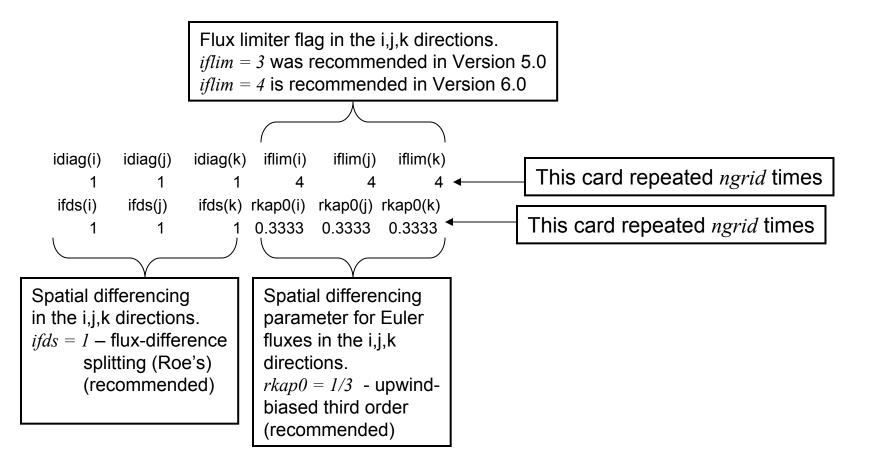


Miscellaneous input



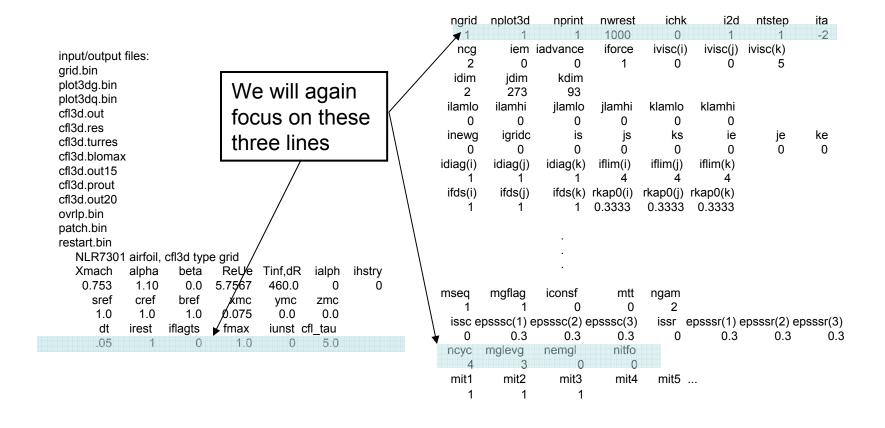


Miscellaneous input



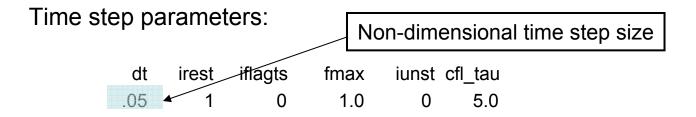


Input for time advancement

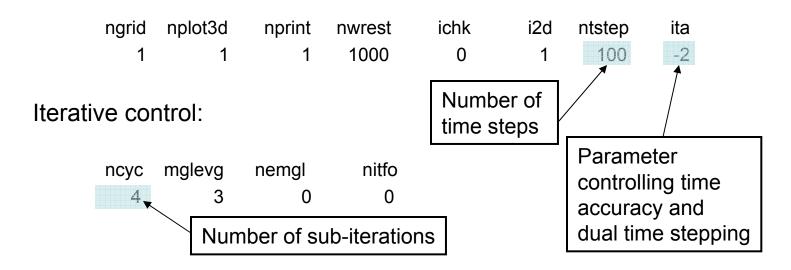




Input for time advancement



Number of time step advances, and time accuracy:





Input for time advancement

Order of time-accuracy, dual time scheme flag (ita)

ita = +1	First order accurate in time; physical time term only
	(t-TS) method
ita = +2	Second order accurate in time; physical time term only
	(t-TS) method
ita = -1	First order accurate in time; physical time and pseudo time term ($\tau\text{-TS}$) method
ita = -2	Second order accurate in time; physical time and pseudo time term (τ-TS) method



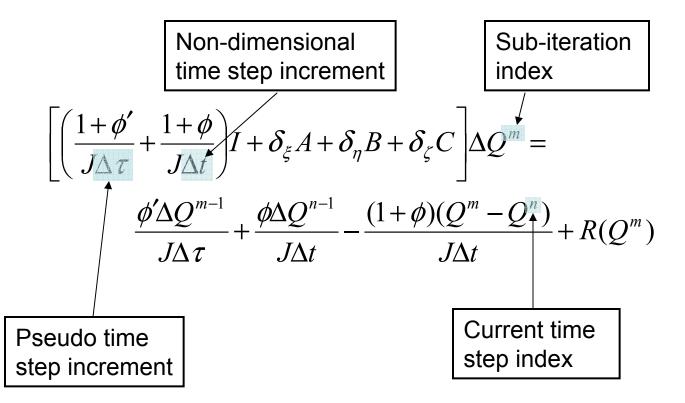
Input for time advancement

Note:

- The approximate factorization scheme used to advance the solution in time introduces first order errors in time. Furthermore, if the diagonal version is utilized (idiag = 1), additional errors of order $\Delta \tau$ are introduced. Sub-iterations can be used to drive these factorization errors to zero. Therefore, if a formally second-order (in time) solution is desired, sub-iterations must be used.
- The inclusion of a pseudo time term increases (often dramatically) the
 maximum allowable time step one can take for a particular problem. However,
 sub-iterations (ncyc > 1) are therefore mandatory and multi-grid is highly
 recommended.
- Larger time steps imply greater error, therefore second order is recommended.
- You will almost never want to use the t-TS method of time stepping.



Equations for τ -TS time advancement





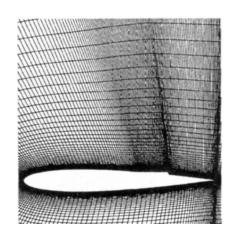
Equations for t-TS time advancement

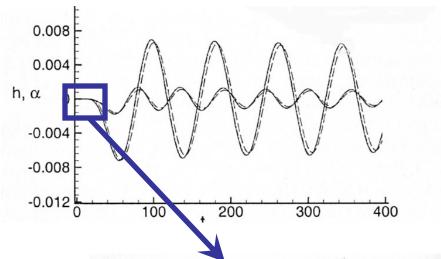
The pseudo time terms are omitted for t-TS time advancement:

$$\begin{bmatrix} \left(\frac{1+\phi}{J\Delta t}\right)I + \delta_{\xi}A + \delta_{\eta}B + \delta_{\zeta}C \end{bmatrix} \Delta Q^{m} = \frac{\phi\Delta Q^{n-1}}{J\Delta t} - \frac{(1+\phi)(Q^{m}-Q^{n})}{J\Delta t} + R(Q^{m})$$
 Non-dimensional time step increment



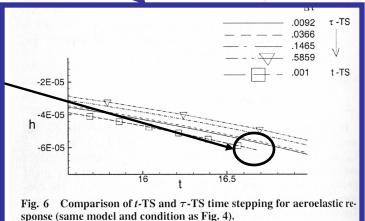
Case study: The t-TS and τ -TS schemes, oscillating spoiler





The solution using the t-TS scheme blows up even at a very small time step size

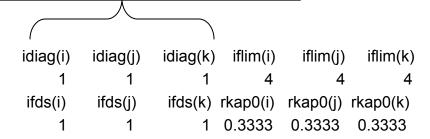
From: Bartels, R. E., "Mesh Strategies for Accurate Computation of Unsteady Spoiler and Aeroelastic Problems," Journal of Aircraft, Vol. 37, No. 3, pp. 521-525.





Speeding up execution time

Parameters controlling the form of the Jacobian matrices used on the left hand side of the equations



Setting idiag(i), idiag(j), idiag(k) to 1 results in a very efficient trigiagonal inversion of the left hand side of the equations in the i, j and k directions. However, be aware of the implications of setting this



Diagonalized versus full Jacobian matrices

idiag controls the form of the matrices A, B, C on the left hand side only. If idiag = 0, the full 5x5 matrix is used. If idiag = 1, the matrix is diagonalized (i.e. Very efficient scalar tridiagonal inversion of the left hand side of this equation).

$$\left[\left(\frac{1 + \phi'}{J\Delta \tau} + \frac{1 + \phi}{J\Delta t} \right) I + \delta_{\xi} A + \delta_{\eta} B + \delta_{\zeta} C \right] \Delta Q^{m} = \frac{\phi' \Delta Q^{m-1}}{J\Delta \tau} + \frac{\phi \Delta Q^{n-1}}{J\Delta t} - \frac{(1 + \phi)(Q^{m} - Q^{n})}{J\Delta t} + R(Q^{m}) \right]$$

Since $\Delta Q^m \to 0$ when the solution converges, setting idiag = 1 does not affect accuracy, ... assuming the solution has been adequately converged.



Sizing *∆t*, number of subiterations

Recall the non-dimensionalization of time:

$$\Delta t = \frac{\Delta \widetilde{t} \ \widetilde{a}_{\infty}}{\widetilde{L}_{R}}$$

The reference length \widetilde{L}_R will be determined by the grid. For instance, if a wing with a 5 inch physical chord length is modeled with a grid that has a non-dimensional chord length of 5, then

$$\widetilde{L}_R = \frac{5 \text{ inches}}{5} = 1 \text{ inch}$$

Note that in this case speed of sound, \tilde{a}_{∞} must be in inches/second.



Sizing *∆t*, number of subiterations

- One criteria for time step sizing is the time scale required to resolve a
 phenomenon at some frequency. Another is the number of time steps
 for a flow field particle to pass over a chord length. Consider 100 time
 steps per cycle or 100 time steps to pass over a chord length as the
 absolute minimum, which ever is smaller.
- The time step size and the number of sub-iterations may have to be set lower/higher respectively by either accuracy or robustness requirements. Short test runs should be performed to ensure adequate convergence.



Sizing *∆t*, number of subiterations

- Indicators that the time step size is too large:
 - The solution converges very slowly or does not converge at all.
 - The solution simply blows up.
 - There are large numbers of negative turbulence parameter values in the file 'cfl3d.subit_turres' the number of which is not converging toward zero at the end of each time step.
- Indicator that the number of sub-iterations is too small:
 - The force coefficients have not leveled out to an acceptable convergence level.
 - The residuals have dropped only by an insufficient magnitude. This can also be a sign that the time step is too large.
 - The solution has been converging, but eventually blows up or starts to gradually diverge.
- Note that these symptoms can also be due to problems with the grid, boundary conditions or turbulence model, so first ensure these issues are settled.



Sub-iterative output – checking convergence

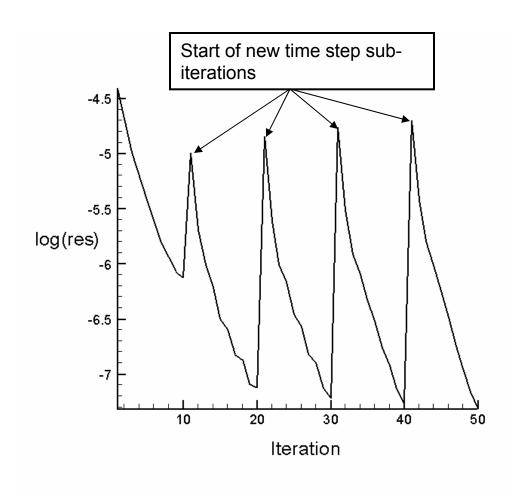
The file 'cfl3d.subit_res' contains the following sub-iterative output

```
subit log(subres)
                            1 -0.44098E+01 -0.56246E-02 0.29632E+00 0.00000E+00 0.14528E-02
                            2 -0.45238E+01 0.28737E-01 -0.12683E-01
                                                                     0.00000E+00 -0.50177E-02
                            3 -0.49884E+01 0.26860E-01 0.19477E+00 0.00000E+00 -0.47901E-02
ncyc = 10 so there
                            4 -0.48541E+01 0.25869E-01 0.80380E-01
                                                                     0.00000E+00 -0.42342E-02
                            5 -0.54203E+01 0.26254E-01 0.10470E+00 0.00000E+00 -0.42906E-02
are 10 lines output
                            6 -0.53829E+01 0.27267E-01 0.98269E-01
                                                                     0.00000E+00 -0.44789E-02
per time step
                            7 -0.58126E+01 0.27020E-01 0.10995E+00 0.00000E+00 -0.44088E-02
                            8 -0.57635E+01 0.26710E-01 0.10469E+00 0.00000E+00 -0.43687E-02
                            9 -0.60754E+01 0.26657E-01 0.10302E+00 0.00000E+00 -0.43724E-02
                           10 -0.61285E+01 0.26713E-01 0.10312E+00 0.00000E+00 -0.43877E-02
                           11 -0.49984E+01 0.26728E-01 0.10431E+00 0.00000E+00 -0.43800E-02
                           12 -0.56927E+01 0.26415E-01 0.92217E-01
                                                                     0.00000E+00 -0.42151E-02
                           13 -0.60126E+01 0.26287E-01 0.83844E-01
                                                                     0.00000E+00 -0.40628E-02
                           14 -0.62182E+01 0.26167E-01 0.82317E-01
                                                                     0.00000E+00 -0.40236E-02
                           15 -0.65022E+01 0.26110E-01 0.82955E-01
                                                                     0.00000E+00 -0.40152E-02
                           16 -0.65972E+01 0.26076E-01 0.83164E-01
                                                                     0.00000E+00 -0.40164E-02
                           17 -0.68247E+01 0.26050E-01 0.82959E-01
                                                                     0.00000E+00 -0.40162E-02
                           18 -0.68719E+01 0.26052E-01 0.82589E-01
                                                                     0.00000E+00 -0.40151E-02
                           19 -0.70916E+01 0.26059E-01 0.82439E-01
                                                                     0.00000E+00 -0.40141E-02
                           20 -0.71274E+01 0.26055E-01 0.82404E-01
                                                                     0.00000E+00 -0.40133E-02
```

Note that all iterations are output sequentially

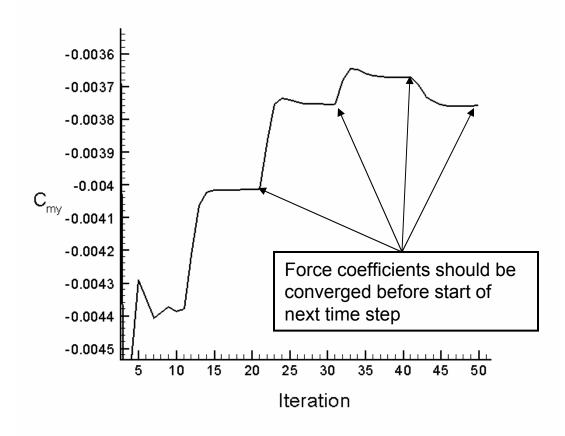


Sub-iterative output—checking convergence





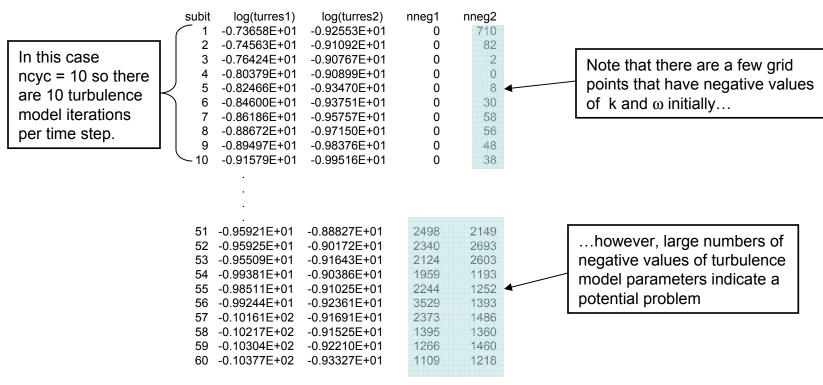
Sub-iterative output—checking convergence





Sub-iterative turbulence output

The file 'cfl3d.subit_turres' contains the following sub-iterative output for Menter's shear stress transport (SST) k-w turbulence model:

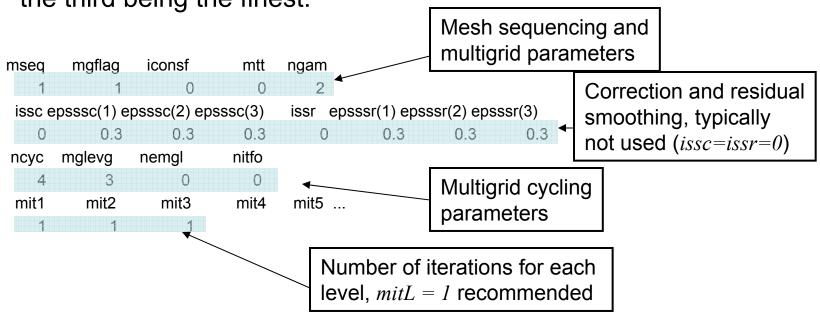


Even though the turbulence model appears to be converging well, a large number of negative values may mean that the time step size is too large for the turbulence model. Usually reducing time step size will fix this problem.



Multigrid strategies

Multigrid is a must for unsteady computations. The following input section establishes four multigrid sub-iterations each on three levels, the third being the finest:





Multigrid strategies

mseq	mgflag	iconsf	mtt	ngam			
1	1	0	0	2			
issc	epsssc(1)	epsssc(2)	epsssc(3)	issr	epsssr(1)	epsssr(2)	epsssr(3)
0	0.3	0.3	0.3	0	0.3	0.3	0.3
ncyc	mglevg	nemgl	nitfo				
4	3	0	0				
mit1	mit2	mit3	mit4	mit5			
1	1	1					

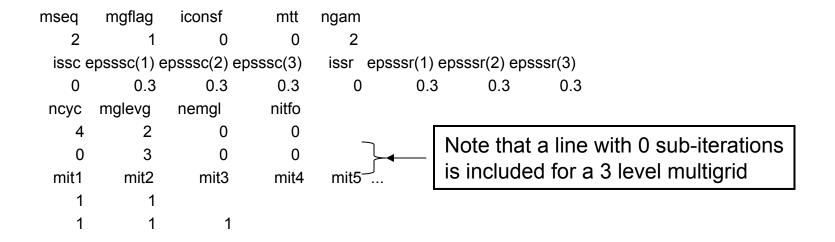
Note:

- *iconsf* is a parameter for setting conservative flux treatment for embedded grids. For most computations it is set to zero.
- *mtt* is a flag for additional iterations on the up portion of the multigrid. Recommend setting to zero.
- ngam is the multigrid cycle flag. ngam = 1 sets V-cycle, ngam = 2 sets a W-cycle. The W-cycle is not recommended for overlapped grids.
- mglevg is the number of grids to use in multigrid cycling. E.g. mglevg = 1 sets the finest grid level only, mglevg = 2 sets two grid levels, etc...
- *nemgl* is set to zero when there are no embedded grids.
- *nitfo1* is the number of first order iterations. Zero is recommended.



Multigrid strategies

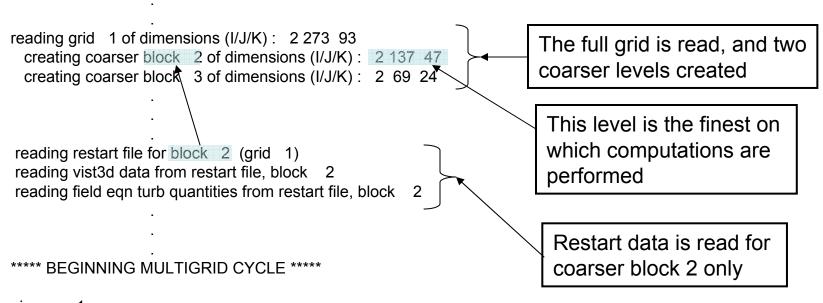
What if you want to compute an unsteady solution using multigrid on coarser levels only? Assume that the steady starting solution has been performed on coarser levels only, as we previously discussed. The following input will allow you to perform the unsteady run:





Multigrid strategies

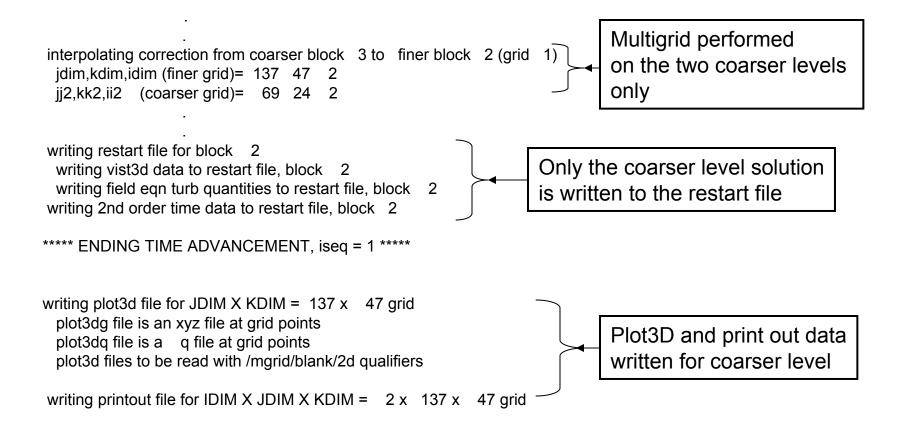
....and here is the output:



iseq= 1 level top = 2 level bottom = 1 number of global grid levels = 2 lglobal= 2



Multigrid strategies



User Specified Grid Motion Overview



CFL3D has the capability to perform computations for prescribed surface motion in two ways

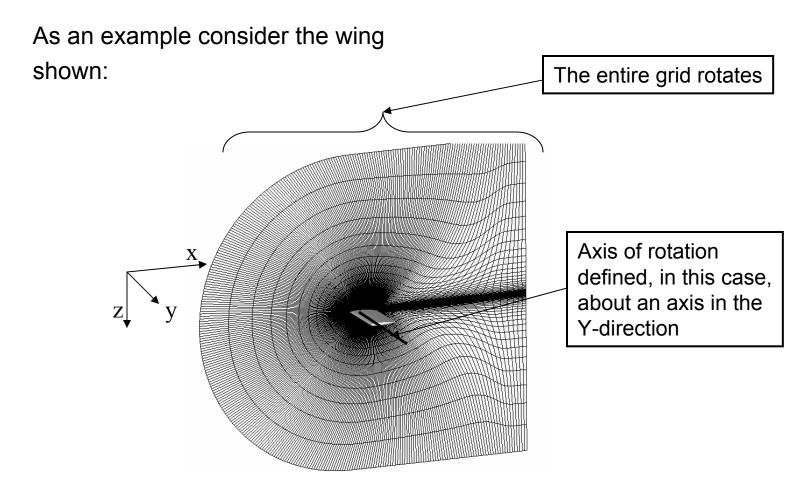
- 1. Prescribed, or user specified rigid grid motion. In this mode, the entire grid or set of grids translates or rotates in a manner prescribed by user input.
- 2. Prescribed surface motion with deforming mesh. In this mode, the surface(s) prescribed by the user translate or rotate and the mesh deforms accordingly.

These types of motion are only available when the code is running in unsteady mode.

User Specified Grid Motion



Rigid grid rotation





Rigid grid rotation

The following unsteady input file performs rotation about the axis shown:

```
input/output files:
wbgrid.cfl
plot3dg.bin
plot3dq.bin
cfl3d.out
cfl3d.res
cfl3d.turres
cfl3d.blomax
cfl3d.out15
cfl3d.prout
cfl3d.out20
ovrlp.bin
patch.bin
restart.bin
NASA Langley BACT Model: NACA 0012 af, AR=1.5 wing, 75TE Flap
             alpha
                                  ReUe
                                           Tinf,dR
                                                              ihstry
                        beta
 0.82000 0.00000 0.00000 0.236E+07
                                           486.00
     sref
               cref
                        bref
                                    xmc
                                              ymc
                                                        zmc
   1.000
          1.00000
                    1.00000
                                0.25000
                                          0.00000
                                                   0.00000
                                             iunst cfl_tau
              irest
                       iflagts
                                   fmax
                                                     2.00000
 0.04000
                       3000
                                 1.00000
    ngrid nplot3d
                      nprint
                                  nwrest
                                              ichk
                                                         i2d
                                                               ntstep
                                                                         ita
                                    1000
                                                0
                                                                         -2
```

Note that iunst = 1 for rigid translation or rotation



Rigid grid rotation

	ivisc(k)	ivisc(j)	ivisc(i)	iforce	iadvance	· ·	
	5	5	5	1	0	2 0	2
					kdim	, ,	idim
					73		73
		klamhi	klamlo	jlamhi	jlamlo	ilamhi	ilamlo
		0	0	0	0	0	0
ke	je	ie	ks	js	is	igridc	inewg
0	0	0	0	0	0	0	0
		iflim(k)	iflim(j)		idiag(k)	idiag(j)	idiag(i)
		3	3	3	1	1	1
			rkap0(j)			fds(j)	ifds(i)
		0.3333		0.3333	1	1	1
iovrlp	nbckdim	nbck0	nbcjdim	nbcj0	nbcidim	nbci0	grid
0	1	5	1	1	1	1	1
ndata	kend	ksta	jend	jsta	bctype	segment	i0: grid
0	73	1	345	1	1001	1	1
ndata	kend	ksta	jend	jsta	bctype	segment	idim: grid
0	73	1	345	1	1002	1	1
ndata	kend	ksta	iend	ista	bctype	segment	j0: grid
0	73	1	73	1	1003	1	1
ndata	kend	ksta	end	ista	bctype	segment	jdim: grid
0	73	1	73	1	1003	1	1
ndata	jend	jsta	iend	ista	bctype	segment	k0: grid
0	33	1	49	1	0	1	1
2	313	33	49	1	2004	2	1
						cq	tw/tinf
						0.00000	0.00000
0	345	313	49	1	0	3	1
0	173	1	73	49	0	4	1
0	345	173	73	49	0	5	1
ndata	jend	jsta	iend	ista	bctype	segment	kdim: grid
0	345	1	73	1	1003	1	1

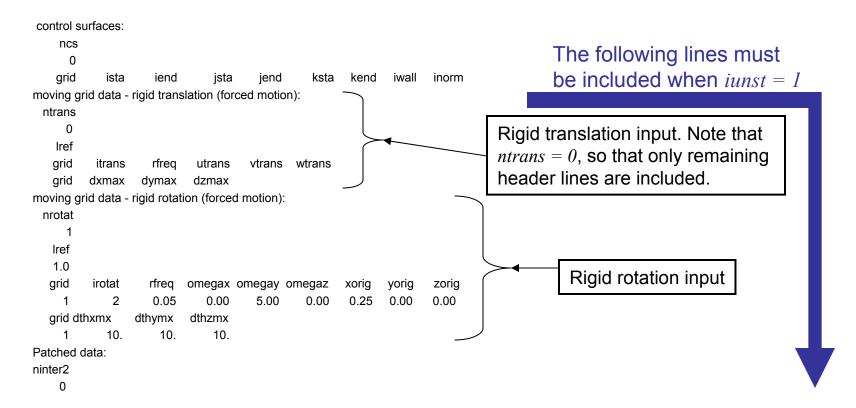


Rigid grid rotation

mseq 1	mgflag 2	iconsf 1	mtt 0	ngam 2						
issc e	nsssc(1) e	epsssc(2)	ensssc(3)	issr	ensssr(1)	epsssr(2)	ensssr(3)			
0	0.3000	0.3000	0.3000	0		0.3000	0.3000			
ncyc	mglevg	nemgl	nitfo	·	0.0000	0.0000	0.0000			
8	3	0	0							
mit1	mit2	mit3	mit4	mit5						
1	1	1	IIIICT	111110						
1-1 block	ing data:									
nbli										
2										
number	grid	ista	jsta	ksta	iend	jend	kend	isva1	isva2	
1	1	1	1	1	49	33	1	1	2	
2	1	49	1	1	73	173	1	1	2	
number	grid	ista	jsta	ksta	iend	jend	kend	isva1	isva2	
1	1	1	345	1	49	313	1	1	2	
2	1	49	345	1	73	173	1	1	2	
patch inte	erface data	a:								
ninter										
0										
plot3d ou	tput:									
grid	iptyp	ista	iend	iinc	jsta	jend	jinc	ksta	kend	kinc
1	0	1	49	1	1	345	1	1	1	1
movie										
0										
print out:										
grid	iptyp	ista	iend	iinc	jsta	jend	jinc	ksta	kend	kinc
⁻ 1	0	1	49	1	1	345	1	1	1	1



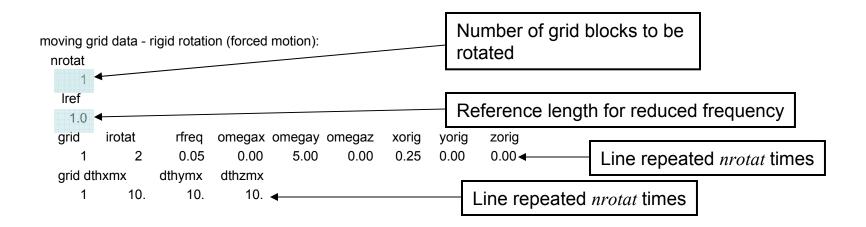
Rigid grid rotation input





Rigid grid rotation input

Focusing attention on the rigid rotation input:





Rigid grid rotation input

Focusing on the last two lines of input on the last slide:

		•						
grid	irotat	rfreq	omegax	omegay	omegaz	xorig	yorig	zorig
1	2	0.05	0.00	5.00	0.00	0.25	0.00	0.00
grid	dthxmx	dthymx	dthzmx					
1	10.	10.	10.					

grid

- Grid block to be rotated

irotat

- Type of rotation

= 0

- no rotation

= 1

- rotation with constant angular speed

= 2

- sinusoidal variation of angular displacement

= 3

- smooth increase in displacement,

asymptotically reaching a maximum angle

rfreq

- reduced frequency when irotat = 2; growth rate to maximum angular displacement when irotat = 3



Rigid grid rotation input

grid irotat rfreq omegax omegay omegaz zorig 0.05 0.00 5.00 0.00 0.00 0.00 grid dthxmx dthymx dthzmx 10. 10. 10.

-

omegax, omegay, omegaz - x,y,z components of rotational velocity when irotat = 1; maximum angular displacements about x,y,z-axes when irotat > 1

xorig, yorig, zorig

- x,y,z coordinate of origin of the rotational axis

dthymx, dthymx, dthzmx

- maximum (absolute) rotational displacement about the x,y,z-axes to be allowed for this grid (set dthxmx,dthymx,dthzmx = 0 if no restriction is required)



Rigid grid rotation input

Example of sinusoidal rotational motion irotat = 2. The following terms are defined:

$$\begin{split} \textit{rfreq} &= k_r \quad , \quad \textit{lref} = L_{\textit{ref}} \\ \textit{omegax} &= \widetilde{\theta}_{x,\text{max}}, \deg. \quad , \quad \textit{omegay} = \widetilde{\theta}_{y,\text{max}}, \deg. \quad , \quad \textit{omegaz} = \widetilde{\theta}_{z,\text{max}}, \deg. \end{split}$$

The rotational displacement (radians) within the code is governed by

$$\theta_{x} = \widetilde{\theta}_{x,\text{max}} \frac{\pi}{180} \sin(2\pi k_{r} \frac{t}{L_{ref}}) , \quad \theta_{y} = \widetilde{\theta}_{y,\text{max}} \frac{\pi}{180} \sin(2\pi k_{r} \frac{t}{L_{ref}})$$

$$\theta_{z} = \widetilde{\theta}_{z,\text{max}} \frac{\pi}{180} \sin(2\pi k_{r} \frac{t}{L_{ref}})$$

Based on these equations of sinusoidal motion,

$$\Delta t = \frac{L_{ref}}{k_r N}$$

where *N* is the desired number of time steps per cycle. Consult Chapter 4 of the Version 5.0 User's Manual pp. 55-62 for details on all types of motion.



Rigid grid rotation

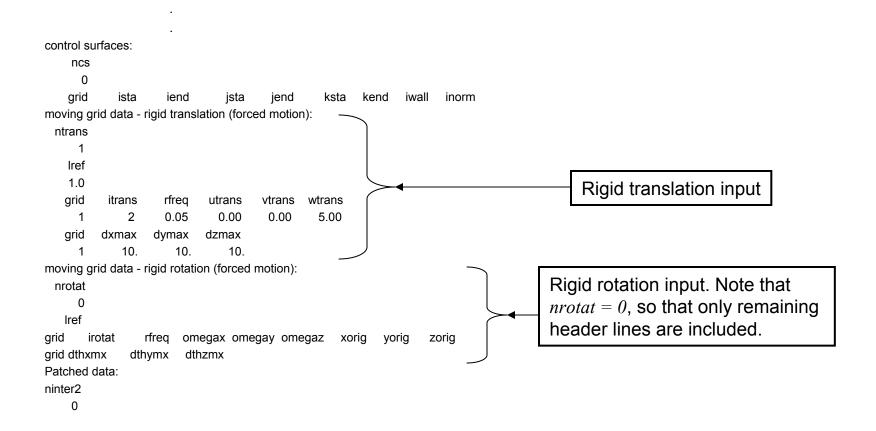
The following diagnostic information on the rotation of the surface(s) will be printed in 'cfl3d.out':

rotating block 1 to new position
creating coarser block 2 of dimensions (I/J/K): 37 173 37
restricting grid speeds from finer block 1 to coarser block 2
creating coarser block 3 of dimensions (I/J/K): 19 87 19
restricting grid speeds from finer block 2 to coarser block 3

Writing restart file for block 1
writing vist3d data to restart file, block 1
writing 2nd order time data to restart file, block 1
writing dynamic mesh data to restart file, block 1
writing dynamic mesh data to restart file, block 1



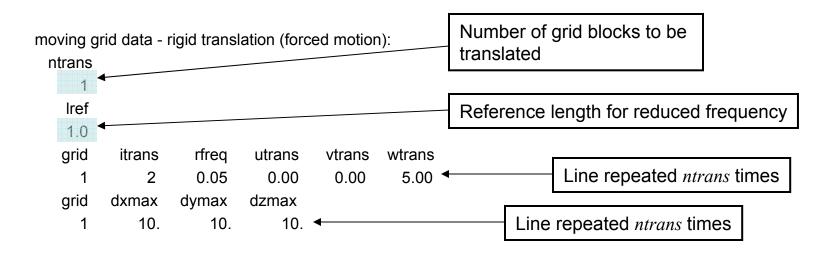
Rigid grid translation input





Rigid grid translation input

Focusing attention on the rigid translation input:





Rigid grid translation input

Focusing on the last two lines of input from the last slide:

grid	itrans	rfreq	utrans	vtrans	wtrans
1	2	0.05	0.00	0.00	5.00
grid	dxmax	dymax	dzmax		
1	10.	10.	10.		

grid Grid block to be rotated

- Type of translation itrans

dymax, dymax,dzmax

= 0no translation

- translation with constant speed = 1

= 2 - sinusoidal variation of displacement

= 3 - smooth increase in displacement, asymptotically reaching a maximum displacement

rfreq - reduced frequency when itrans = 2; growth rate to maximum displacement when itrans = 3

utrans, vtrans, wtrans

- x,y,z components of translation velocity when *itrans* = 1; maximum

displacements in the x,y,z directions when itrans > 1

- maximum (absolute) translation displacement in the x,y,z directions to be allowed for this grid.



Overview

- CFL3D can perform several types of user specified surface motion by deforming the mesh, i.e. surface rotation and/or translation of all or partial segments of the solid surfaces as well as modal motion of surfaces.
- Aeroelastic, user defined deforming mesh surface and user defined rigid grid motion can be performed in any combination.
- There are two methods of deforming the mesh.
 - Exponential decay combined with Trans-Finite Interpolation (TFI) of interior mesh points.
 - Finite macro-element deformation combined with TFI.
- Note that deforming surface motion can only be performed with the code running in unsteady mode.

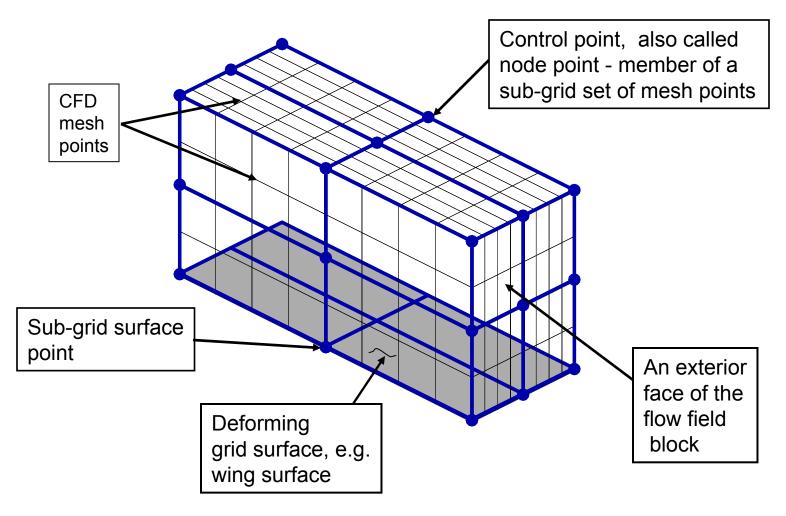


Overview

- In the first mesh movement option (exponential decay method) deformation is performed in two steps.
 - The first step is exponential decay of control points away from the moving surface. The rate
 of the exponential decay is controlled by user input.
 - The second step is a TFI of mesh points interior to the control points.
- Advantage of the exponential decay method is that it is computationally efficient
- In the second mesh movement option (finite macro-element method) deformation is also performed in two steps.
 - The first step is a finite element solution of macro-element points. The resulting solution transmits surface motion to the element node points. The element stiffness varies with distance from the surface. User specified input controls the rate at which the element stiffness decays away from surfaces.
 - The second step is a TFI of mesh points interior to the element node (or control) points.
 - See Bartels, R. E., "Finite Macro-Element Mesh Deformation in a Structured Multi-Block Navier-Stokes Code," NASA/TM-2005-213789, July 2005.
- Advantage of the finite macro-element method is that it maintains mesh quality, but is significantly more computationally time consuming.

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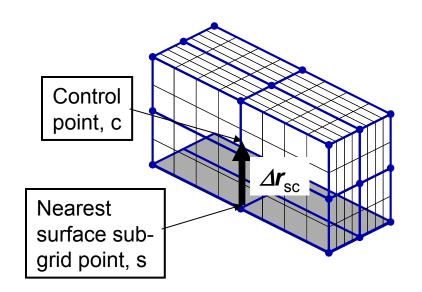
Deforming mesh terminology





Deforming mesh using exponential decay method

$$\vec{r}_c^{n+1} - \vec{r}_c^n = D_{sc}(\vec{r}_s^{n+1} - \vec{r}_s^n)$$
where
$$D_{sc} = \min[1, e^{-A}]$$
and
$$A = \beta_2 \left(\left| \Delta \vec{r}_{sc} \right| / \Delta r_{\text{max}} - \alpha_2 \right)$$



The movement of surface points is transmitted into the flow field sub-grid through an exponential decay function D_{sc} . The rate of decay is controlled by the parameters β_2 and α_2 .

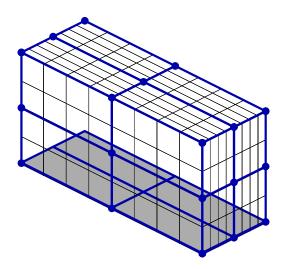


Deforming mesh with exponential decay method

Note several potential draw backs to this approach:

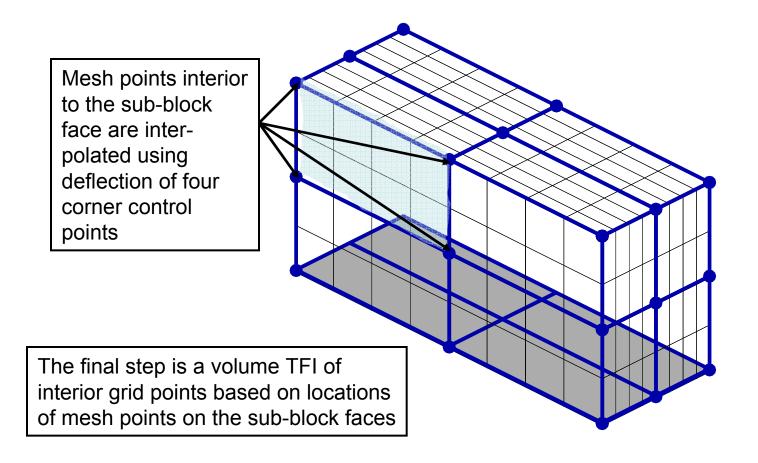
- Too rapid a rate of decay (β_2 too large, α_2 too small) results in the possibility of the surface points moving through nearby control points.
- Too low a rate of decay (β_2 too small, α_2 too large) results in the possibility of surface deformation being transmitted too far into the flow field with possible penetration of opposing surfaces.
- Typical values for decay parameters are:

$$\beta_2 = 1 - 10$$
 , $\alpha_2 = 0.005 - 0.05$



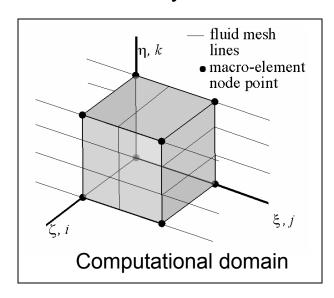


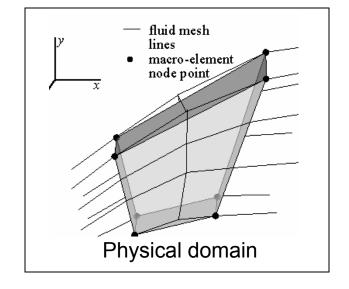


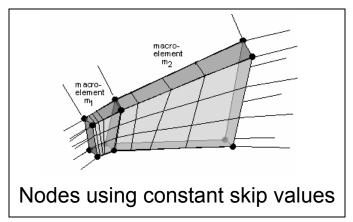


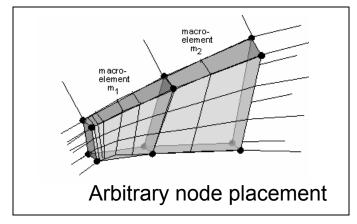


Coordinate systems and terminology for finite macro-element method









Finite macro-element method

The equations of elasticity are solved using Hooke's law for element m

$$\vec{\sigma}_m = C_m \vec{\varepsilon}_m$$

where

$$ec{\sigma}_{m} = \begin{bmatrix} \sigma_{xx} \\ \sigma_{yy} \\ \sigma_{zz} \\ \sigma_{xy} \\ \sigma_{yz} \end{bmatrix}_{m}, \quad \vec{\mathcal{E}}_{m} = \begin{bmatrix} \mathcal{E}_{xx} \\ \mathcal{E}_{yy} \\ \mathcal{E}_{zz} \\ \mathcal{E}_{xy} \\ \mathcal{E}_{xz} \\ m \end{bmatrix}_{m}, \quad C_{m} = \begin{bmatrix} E_{m} & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & E_{m} & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & E_{m} & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & G_{m} & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & G_{m} & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & G_{m} \end{bmatrix}$$

$$E_m = E_0 f_m \quad , \quad G_m = G_0 f_m$$

$$E_{m} = E_{0} f_{m}$$
 , $G_{m} = G_{0} f_{m}$
$$f_{m} = \frac{1}{1 - \exp(-\beta_{1} \Delta r_{m} / \Delta r_{max})}$$

 Δr_m is computed as

$$\Delta r_m = \sqrt{(\Delta x_{cs,m})^2 + (\Delta y_{cs,m})^2 + (\Delta z_{cs,m})^2}$$

The user controls the rate of decay of material properties by the parameter β_1 . Typical values of β_1 are in the range of 1-2.



Input for deforming mesh

Moving grid data – data for field/multiblock mesh movement beta2 nskip isktyp beta1 alpha1 alpha2 nsprgit 4 -1 2.0 1.1 10.0 0.01 0 jskip grid iskip kskip Moving grid data - multi-motion coupling ncoupl 0 Slave master xorig yorig zorig - number of blocks for which skip value data is input. If nskip = 0 the code nskip computes default skip values (isktyp = -1, 1) or control point index values (isktyp = -2,2).- Parameter defining the mesh deformation approach isktyp

exponential decay method

finite macro-element method



Input for deforming mesh

Moving grid data – data for field/multiblock mesh movement nskip isktyp beta1 alpha1 beta2 alpha2 nsprgit -1 2.0 1.0 10.0 0.01 1 0 iskip grid iskip kskip Moving grid data – multi-motion coupling ncoupl 0 Slave master xorig yorig zorig

nsprgit

beta1 - Parameter controlling macro-element stiffness decay (typically 1.0-2.0)

alpha1 - Relaxation parameter for Gauss-Seidel solver (typically 0.8-1.2).

beta2 - Decay parameter for the exponential decay method (typically 1 - 10).

alpha2 - Decay parameter for the exponential decay method (typically 0.005-0.05).

- Number of spring analogy smoothing steps performed with the exponential decay method. This step applies *nsprgit* spring analogy steps to the control points after application of the exponential decay step (typically 0-2).

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Input for deforming mesh

- There are 4 options for the construction of control points.
 - Option 1: Code generated minimum number of control points.

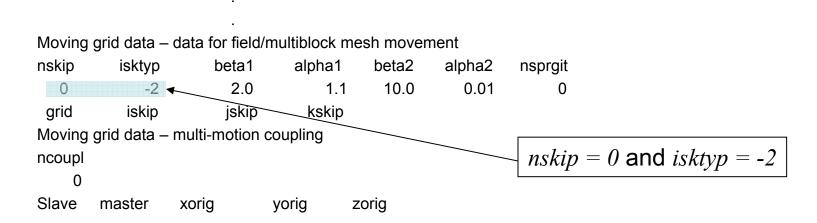
preferred method

- Option 2: Code generated default skip values.
- Option 3: User input of i,j,k skip values for each block.
- Option 4: User defined input of control point i,j,k indices for each block.
- These options depend on the value of *nskip* and the value of *isktyp*
 - Option 1: isktyp = -2, 2 and nskip = 0
 - Option 2: isktyp = -1, 1 and nskip = 0
 - Option 3: isktyp = -1, 1 and nskip = ngrid (Note: ngrid = number of grid blocks)
 - Option 4: isktyp = -2, 2 and nskip = ngrid
- Option 1 creates the minimum number of control points (at non-constant intervals) by
 placing control point points only at each boundary segment extremity. This is the
 preferred method.
- Options 2 creates skip values that result in control points at constant intervals through out each of the grids, with control points at each boundary segment extremity.
 Sometimes this is more robust than option 1, but can create many more control points.



Option 1 – Code generated minimum number of control points

It is possible to have the code calculate the minimum number of control points. This is the preferred method. The following lines of input accomplish that:

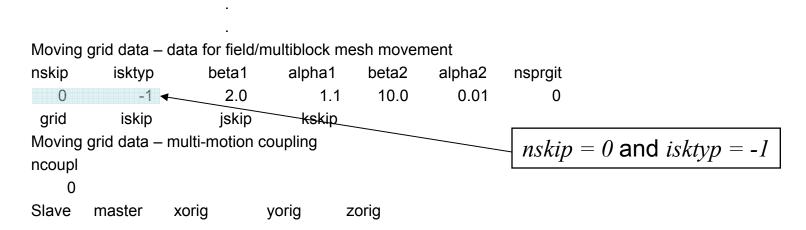


Note that the data input line following the header 'grid' is omitted. The code calculates the minimum number of control points possible consistent with placing control points at each boundary segment extremity. The values it calculates will be found in the 'cfl3d.out' section that reflects input. Note that the value of *isktyp* must be either 2 or -2. In general control points will not be at constant intervals.



Option 2 – Code generated skip values

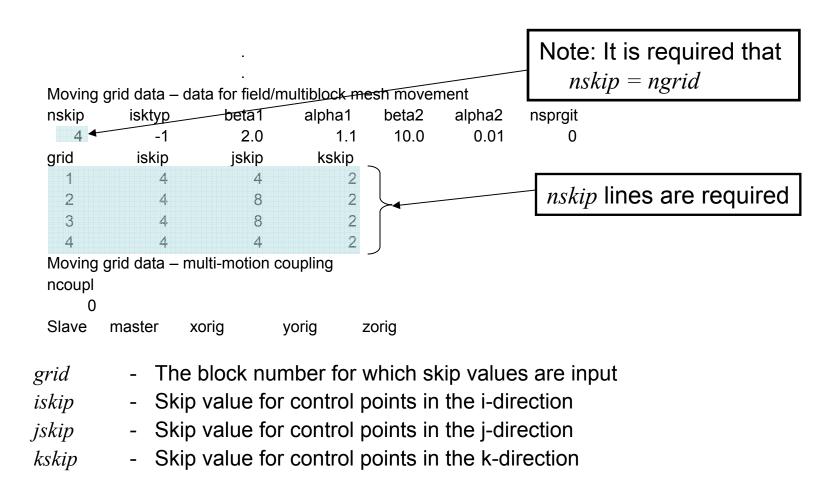
It is possible to have the code calculate default skip values. The following lines of input accomplish that:



Note that the data input line following the header 'grid' is omitted. The code calculates the largest values of *iskip*, *jskip*, *kskip* possible. The values it calculates will be found in the 'cfl3d.out' section that reflects input. Note that the value of *isktyp* must be either 1 or -1.



Option 3 – User *i,j,k* skip input



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Permissible skip values

iskip, jskip, kskip values determine the i, j, k skip intervals for creating the sub-grid

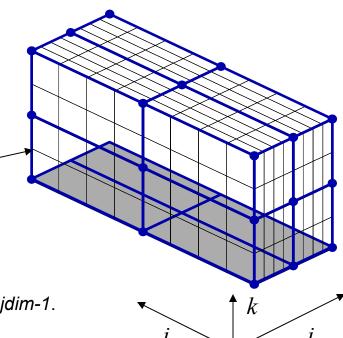
For this grid:

$$idim = 9$$
, $jdim = 9$, $kdim = 5$ and

$$iskip = 4$$
, $jskip = 4$, $kskip = 2$

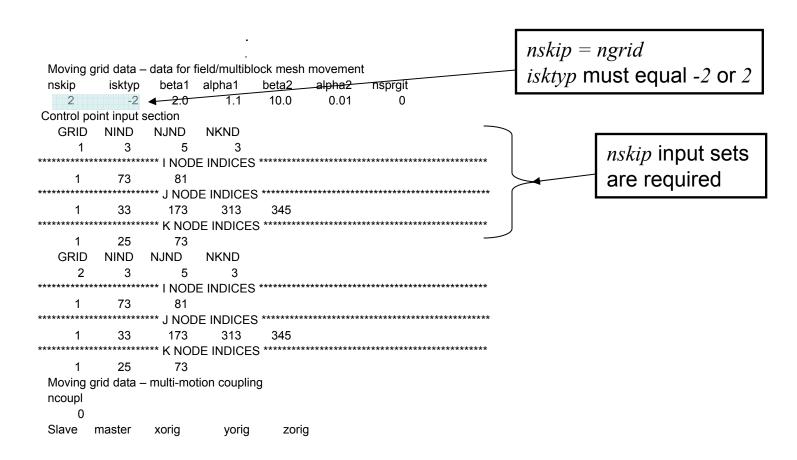
Skip values must evenly divide into one minus the dimension of the grid. *jskip* must divide evenly into *jdim-1*. *iskip* must divide evenly into *idim-1*, etc...

With *idim* = 9, permissible values of *iskip* are 2, 4 and 8. With *jdim* = 9, permissible values of *jskip* are 2, 4 and 8. With *kdim* = 5, permissible values of *kskip* are 2 and 4.





Option 4 – User input of i,j,k control point indices





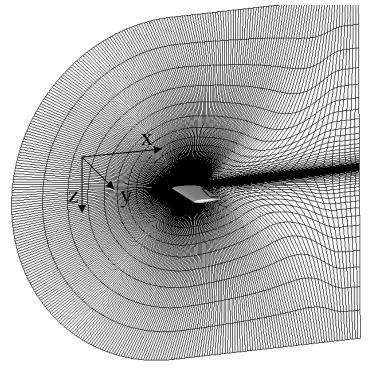
Option 4 — User input of i,j,k control point indices

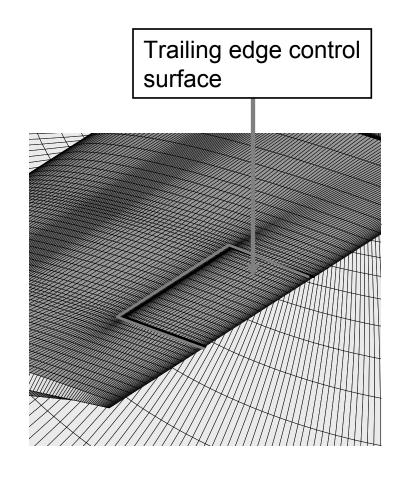
- This option is used when there are problem areas in the surface motion that
 require customized control point placement. e.g. significant surface motion
 restricted to a small portion of the entire surface area or if the finite macro-element
 method is used and added control points are needed to define affine element
 shapes.
- Note that a control point must be placed at the extremities of all boundary condition segments, 1-1 blocking segments and all block corners.
- The code will do a check at 1-1 blocking segments to see if the control points you have selected result in continuity in control placement between 1-1 blocking boundaries. It will add points as necessary to maintain control point continuity. This is a very powerful feature that can be very useful when adding control points.
- The code will not tell you if a b.c. segment extremity or block corner does not have a control point assigned to it. It will simply cause the grid motion to be messed up and produce negative volumes!

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Example 1: 3D Control surface rotation with exponential decay method

As an example consider the wing shown undergoing control surface rotation:







Example 1: 3D Control surface rotation with exponential decay method

The following unsteady input file performs the control surface rotation about the hinge point:

```
input/output files:
 wbgrid.cfl
 plot3dg.bin
 plot3dq.bin
 cfl3d.out
 cfl3d.res
 cfl3d.turres
 cfl3d.blomax
 cfl3d.out15
 cfl3d.prout
 cfl3d.out20
 ovrlp.bin
 patch.bin
 restart.bin
NASA Langley BACT Model: NACA 0012 af, AR=1.5 wing, 75TE Flap
    Mach
                                   ReUe
                                            Tinf,dR
             alpha
                        beta
                                                        ialph
                                                               ihstry
 0.82000 0.00000 0.00000 0.236E+07
                                            486.00
                                                           1
                                                                   0
     sref
               cref
                        bref
                                    xmc
                                              vmc
                                                        zmc
                                           0.00000 0.00000
    1.000 1.00000 1.00000
                                 0.25000
                                                     cfl_tau
       dt
              irest
                       iflagts
                                   fmax
                                              iunst
                                                     2.00000
 0.04000
                       3000
                                 1.00000
    ngrid nplot3d
                       nprint
                                               ichk
                                                          i2d
                                  nwrest
                                                               ntstep
                                                                          ita
                                    1000
                                                 0
                                                                          -2
```

Note that iunst = 2 for deforming mesh

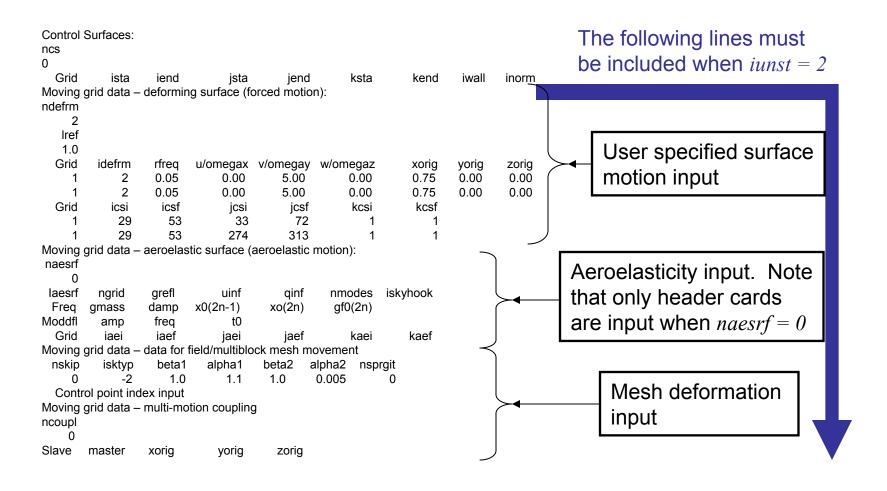


	ncg		iadvance	iforce	٠,	ivisc(j)		
	2		0	1	5	5	5	
	idim	, ,	kdim					
	81	345	73					
i	ilamlo	ilamhi	jlamlo	jlamhi	klamlo	klamhi		
	0	0	0	0	0	0		
i	inewg	igridc	is	js	ks	ie	je	ke
	0	0	0	0	0	0	0	0
id	liag(i)	idiag(j)	idiag(k)	iflim(i)	iflim(j)	iflim(k)		
	1	1	1	4	4	4		
if	fds(i)	fds(j)	ifds(k)	rkap0(i)	rkap0(j)	rkap0(k)		
	1	1	1	0.3333	0.3333	0.3333		
	grid	nbci0	nbcidim	nbcj0	nbcjdim	nbck0	nbckdim	iovrlp
	1	1	1	1	1	5	1	0
i0:	grid	segment	bctype	jsta	jend	ksta	kend	ndata
	1	1	1005	1	345	1	73	0
idim:	grid	segment	bctype	jsta	jend	ksta	kend	ndata
	1	1	1002	1	345	1	73	0
j0:	grid	segment	bctype	ista	iend	ksta	kend	ndata
	1	1	1003	1	81	1	73	0
jdim:	grid	segment	bctype	ista	end	ksta	kend	ndata
•	1	1	1003	1	81	1	73	0
k0:	grid	segment	bctype	ista	iend	jsta	jend	ndata
	1	1	0	1	73	1	33	0
	1	2	2004	1	73	33	313	2
t	w/tinf	cq						
0.0	00000	0.00000						
	1	3	0	1	73	313	345	0
	1	4	0	73	81	1	173	0
	1	5	0	73	81	173	345	0
kdim:	grid	segment	bctype	ista	iend	jsta	jend	ndata
	1	1	1003	1	81	1	345	0



mseq 1 issc e 0 ncyc 8 mit1	mgflag 2 psssc(1) e 0.3000 mglevg 3 mit2	iconsf 1 epsssc(2) 0.3000 nemgl 0 mit3	mtt 0 epsssc(3) 0.3000 nitfo 0 mit4	ngam 2 issr 6 0 mit5	0.3000	epsssr(2) 0.3000	epsssr(3) 0.3000			
1-1 block	ing data:									
nbli 2										
number	grid	ista	jsta	ksta	iend	jend	kend	isva1	isva2	
1	1	1	1	1	73	33	1	1	2	
2	1	73	1	1	81	173	1	1	2	
number	grid	ista	jsta	ksta	iend	jend	kend	isva1	isva2	
1	1	1	345	1	73	313	1	1	2	
2	1	73	345	1	81	173	1	1	2	
patch inter ninter 0	erface data	a:								
plot3d ou	ıtput:									
grid	iptyp	ista	iend	iinc	jsta	jend	jinc	ksta	kend	kinc
1	0	1	73	1	33	313	1	1	1	1
movie 0										
print out:										
grid 1	iptyp 0	ista 1	iend 73	iinc 1	jsta 33	jend 313	jinc 1	ksta 1	kend 1	kinc 1







Example 1: 3D Control surface rotation with exponential decay method

ndefrm 2	ırid data –	deformir	: ng surface (fo	orced motion	n): surfa	ace is a lowe	d
Iref							
1.0							
Grid	idefrm	rfreq	u/omegax	v/omegay	w/omegaz	xorig	

Note that ndefrm = 2 because the trailing edge control surface is defined by an upper wing surface segment and a lower wing surface segment

1.0								
Grid	idefrm	rfreq	u/omegax	v/omegay	w/omegaz	xorig	yorig	zorig
1	2	0.05	0.00	5.00	0.00	0.75	0.00	0.00 ndefrm lines required
1	2	0.05	0.00	5.00	0.00	0.75	0.00	0.00
Grid	icsi	icsf	jcsi	jcsf	kcsi	kcsf		
1	29	53	33	72	1	1	\neg	
1	29	53	274	313	1	1		nacjim iiiles required

Grid

- grid block containing the moving surface

idefrm

- type of surface motion

= 1

- translation

= 2

- rotation

rfreq

- reduced frequency of the surface motion

u/omegax, v/omegay, w/omegaz

- x,y,z-components of surface translational velocity if *idefrm* = 1

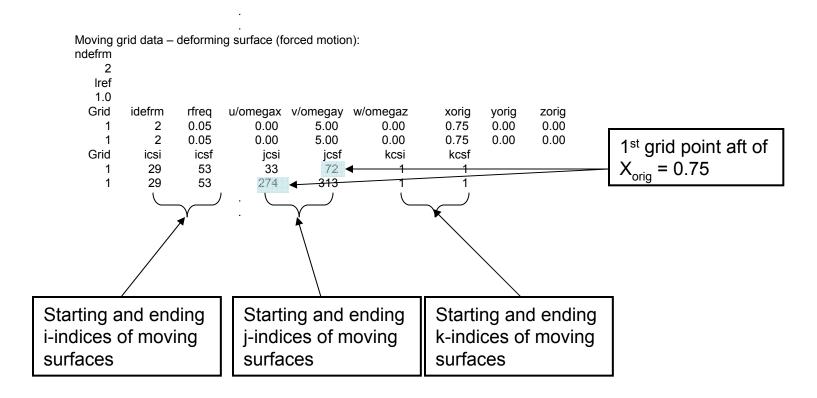
- x,y,z-components of surface rotational velocity if *idefrm* = 2

xorig, yorig, zorig

- x,y,z coordinates of the origin of the rotation axis (note: value must be input even when idefrm = 1)

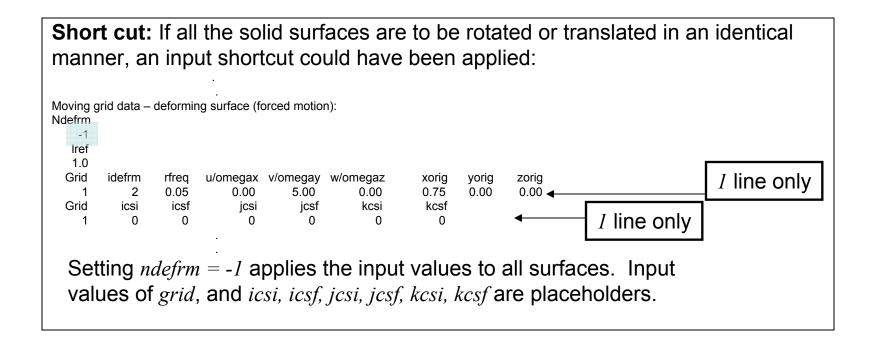


Example 1: 3D Control surface rotation with exponential decay method

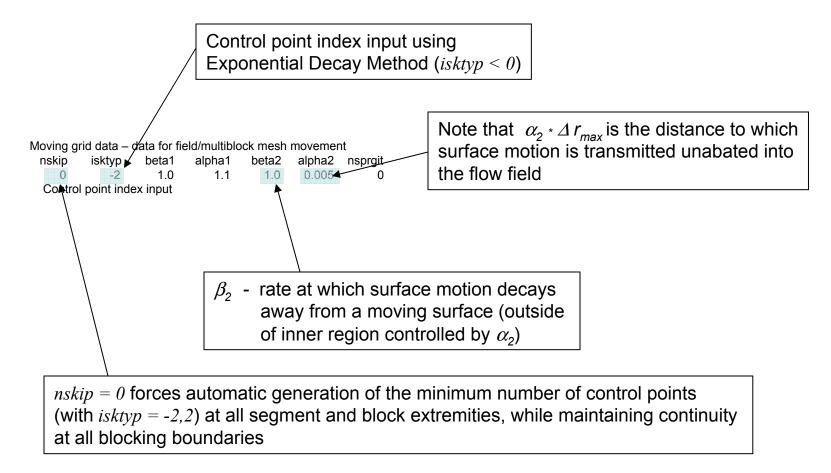


Note that the two surface definitions actually comprise a single control device (upper and lower surfaces of the trailing edge control device).



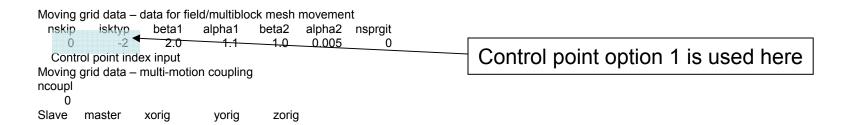








Example 1: 3D Control surface rotation with exponential decay method

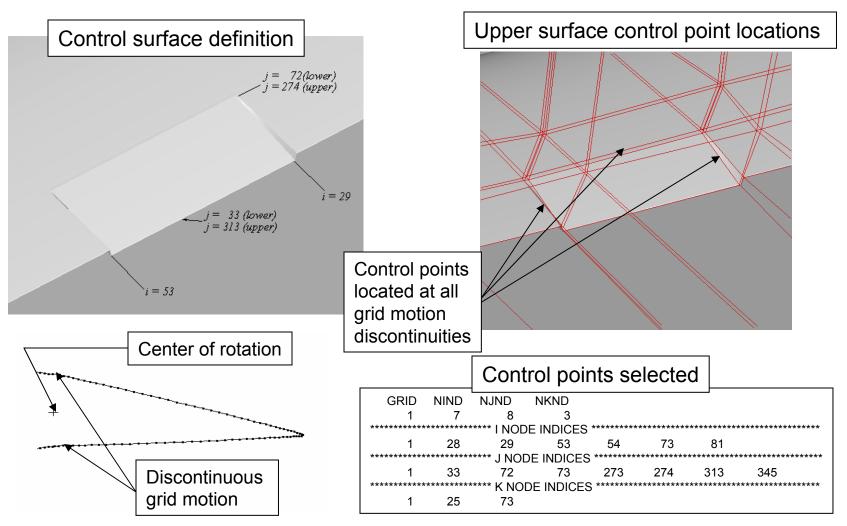


• This input option automatically creates the following control points: (This format is how it would look if you were to input these control points by hand (i.e. using Option 4))

GRID	NIND	NJND	NKND					
1	7	8	2					
*******	******	***** I NOE	E INDICES	******	******	******	******	****
1	28	29	53	54	73	81		
*******	******	***** J NOI	DE INDICES	3 *******	*****	******	******	****
1	33	72	73	273	274	313	345	
*******	******	***** K NO	DE INDICES	S *******	*******	*******	******	****
1	73							

- Note that i node indices, j node indices, k node indices span the entire block. (i.e. idim = 81, jdim = 345, kdim = 73)
- Boundary segments have a control point. The trailing edge at j = 33 and 313 has control points assigned. The wing tip at i = 73 has a control point assigned.
- Other control points have been assigned at discontinuities in the surface movement. (e.g. at i = 28, 29 and 53, 54 and j = 72, 73 and 273, 274) See the next slide.

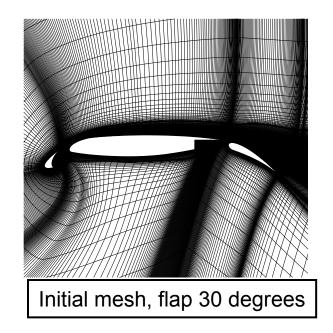


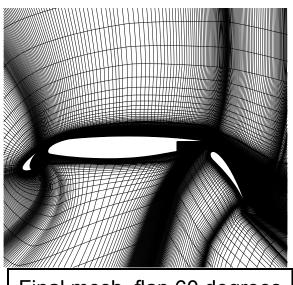




Example 2: 2D Flap rotation with finite macro-element method

Consider the 2D three element airfoil with rotation and translation of the trailing edge flap.



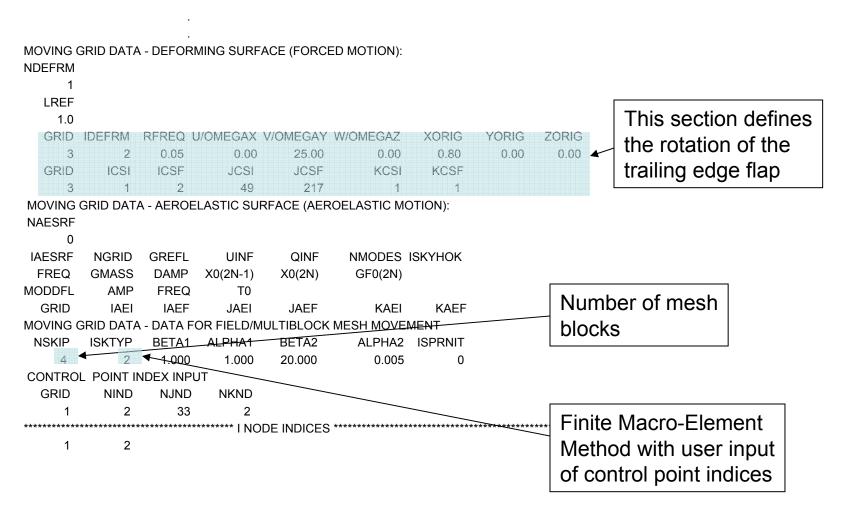


Final mesh, flap 60 degrees

From Bartels, R. E., "Finite Macro-Element Mesh Deformation in a Structured Multi-Block Navier-Stokes Code," NASA/TM-2005-213789, July 2005.



Example 2: 2D Flap rotation with finite macro-element method





Example 2: 2D Flap rotation with finite macro-element method

******	*****	******	*****	J NODE IND	DICES *****	*****	*****	*****	*****
1	10	34	49	75	101	113	137	161	201
237	273	299	317	333	349	380	395	410	433
445	473	509	545	585	609	633	645	671	697
712	736								
*******	******	******	******	K NODE IN	DICES ****	******	******	******	*****
1	57								
GRID	NIND	NJND	NKND						
2	2	27	2						
		******	******	I NODE IND	DICES *****	*****	*****	******	*****
1	2	L-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1	. 	LNODEIN	DIOE0 ****	* * * * * * * * * * * * * * * * * * *		· · · · · · · · · · · · · · · · · · ·	****
				J NODE INI					
1	10		49	75		113		145	157
185	225	261	281	299	325	361	397	437	461
485	497		549		588				
		******	******	K NODE IN	DICES ****	******	*****	*****	*****
1	89								
GRID	NIND		NKND						
3	2	16	2						
		******	******	I NODE IND	DICES *****	*****	*****	*****	*****
1	2								
				J NODE IN					
1	10	34	49	75	101	116	121	129	153
165	191		232		265				
		******	******	K NODE IN	DICES ****	******	*****	*****	*****
1	65								
			NKND						
4	2	32	5						

Up to 10 per line, 500 total allowed

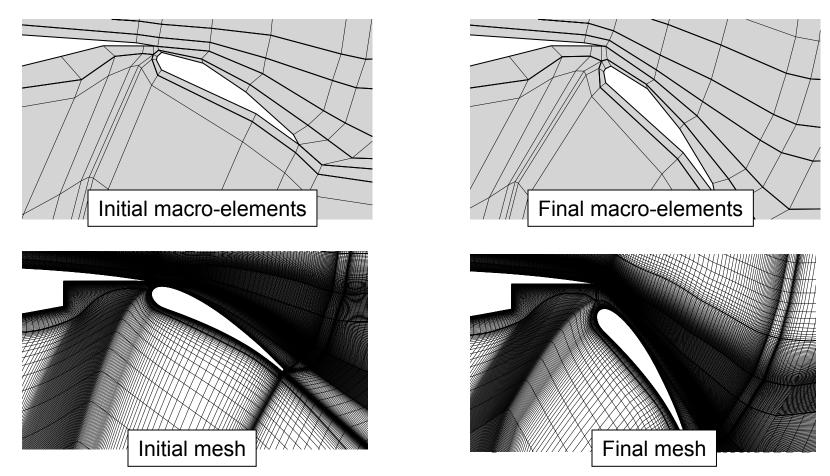


Example 2: 2D Flap rotation with finite macro-element method

*******	*******	*****	*******	NODE IND	ICES *****	*****	******	*****	****
1	2								
*******	******	*****	***** J	NODE IND	DICES *****	******	******	******	****
1	10	34	49	75	101	116	121	133	161
201	237	257	273	289	320	335	350	373	385
413	449	485	525	549	573	585	611	637	652
676	685								
*******	******	******	****** K	NODE IN	DICES ****	*****	******	*****	*****
1	10	17	24	33					
MOVING (GRID DAT	ΓA - MULT	I-MOTION	COUPLIN	NG				
NCOUPL									
0									
SLAVE M	1ASTER	XORIG Y	ORIG ZO	RIG					



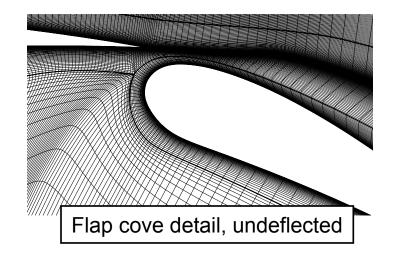
Example 2: 2D Flap rotation with finite macro-element method

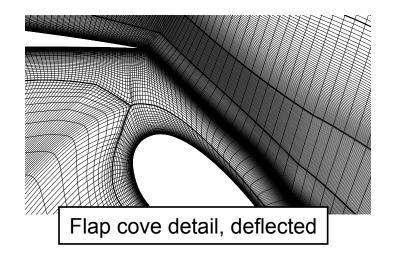


From Bartels, R. E., "Finite Macro-Element Mesh Deformation in a Structured Multi-Block Navier-Stokes Code," NASA/TM-2005-213789, July 2005.



Example 2: 2D Flap rotation with finite macro-element method

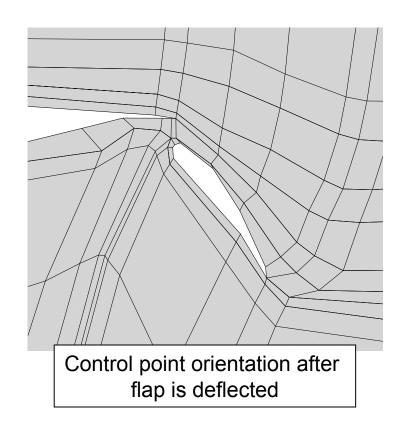


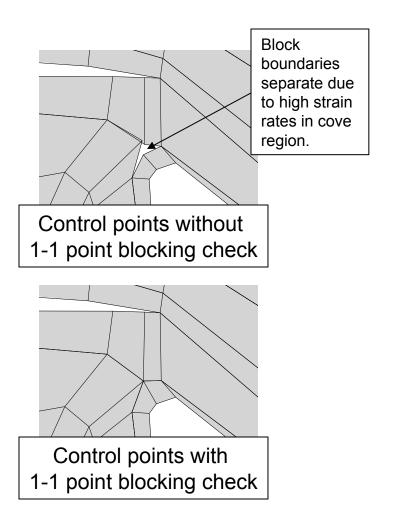


From Bartels, R. E., "Finite Macro-Element Mesh Deformation in a Structured Multi-Block Navier-Stokes Code," NASA/TM-2005-213789, July 2005.



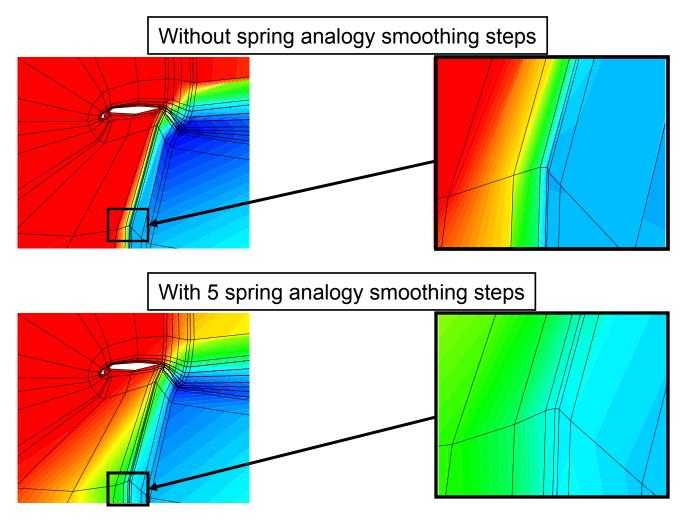
Example 2: 2D Flap rotation – 1-1 block point checking







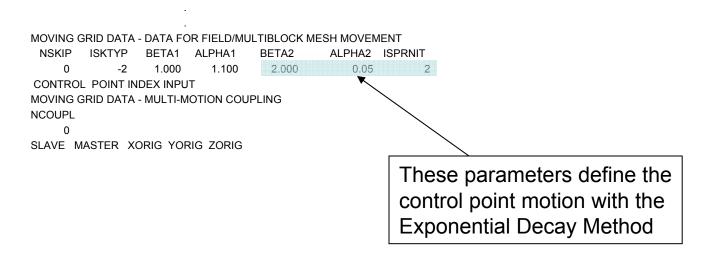
Example 2: 2D Flap rotation using exponential decay method





Example 2: 2D Flap rotation using exponential decay method

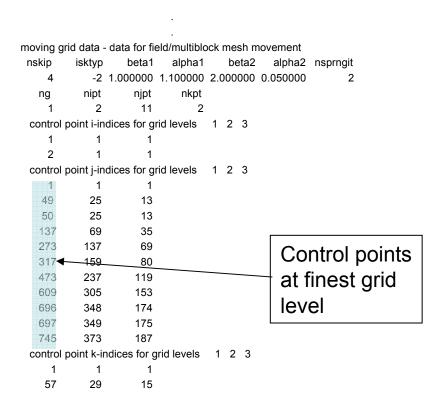
An alternate approach is to allow automatic creation of the minimum number of control points. (Option 1) The input below accomplishes that by setting nskip = 0. Note that the exponential decay method is used (isktyp < 0).





Example 2: 2D Flap rotation using exponential decay method

The control points that are code selected appear in the 'cfl3d.out' file:

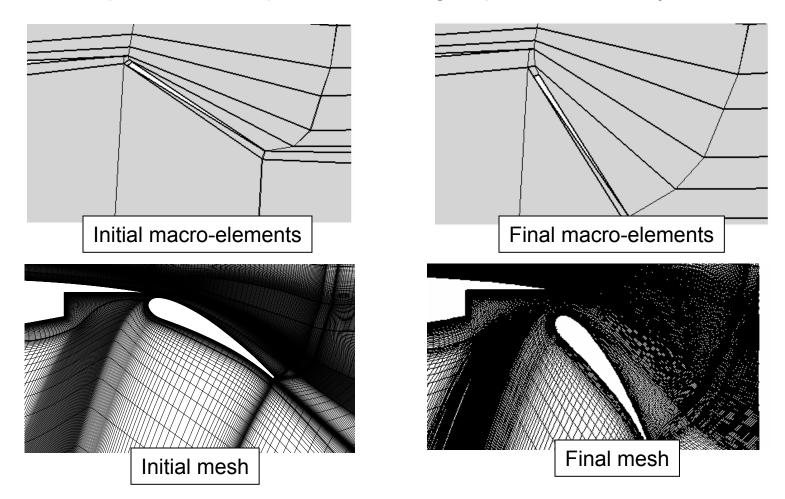


ng	nipt	njpt	nkpt			
2	2	11	2			
control	point i-indic	ces for grid I	evels	4	5	6
1	1	1				
2	1	1				
control	point j-indid	ces for grid I	evels	4	5	6
1	1	1				
49	25	13				
50	25	13				
137	69	35				
145	73	37				
281	141	71				
325	163	82				
461	231	116				
548	274	137				
549	275	138				
597	299	150				
control	point k-indi	ces for grid	levels	4	5	6
1	1	1				
89	45	23				

The resulting mesh movement is shown in the next slide.



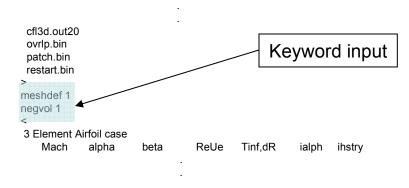
Example 2: 2D Flap rotation using exponential decay method



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Example 2: 2D Flap rotation

- The mesh movement shown in the previous slides is robust (no negative volumes) through the entire range of motion shown, however mesh quality aft of the flap is somewhat degraded after deflection.
- If β_2 is set to 1.0 or if the finite macro-element method is used with the code selected minimum number of control points (as was shown), negative volumes are the result.
- There is a simple way to fix this problem, demonstrated next. In the process an option
 for running the code will be demonstrated in which only the mesh motion and mesh calculations
 (e.g. metric and volume calculations) are performed in the code. This option greatly speeds up the
 code when the mesh motion is being debugged.
- The 'Mesh only' run option is invoked by using the keyword input, *meshdef 1*. Keyword input will be discussed in detail later in the course. Note spelling and capitalization are important.
- The input to accomplish this is as follows:



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Example 2 : 2D Flap rotation

- Setting the keyword *meshdef* to *1* also causes the control points to be output in a Tecplot file in point wise data format. Other auxiliary data are also printed out in other files.
- If one processor is used all block control points are output into the file Tecplot data file 'fort.4000'. Data included in this file are x,y,z locations of control points, x,y,z deflections per time step, node number, and node number of the nearest surface point.
- If multiple processors are used, the control points from the blocks processed on each processor are put in the successive files 'fort.4001, fort.4002, ...'
- Note that if the option movie = inc is used, the control points at every inc time steps will be output. If movie = 0, only control points at the final time step will be output.
- Once the control points are plotted it is possible to better visualize where added control
 points need to be placed.
- This is the option that was used to create the plots of control points shown in this
 presentation.

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Example 2 : 2D Flap rotation

- Returning to the flap rotation example above, say we want to run it using control point option 1 (nskip = 0, isktyp = -2,2) but now using the finite macro-element method (isktyp = 2)
- The input parameters used are: $\beta_1 = 1.0$, $\alpha_1 = 0.9$.
- Keywords 'meshdef 1' and 'negvol 1' are set. When the keyword 'negvol 1' is used, the code continues executing and prints a diagnostic message in 'cfl3d.out' indicating where the negative volume occurred.
- The code encounters negative volumes, with the following messages appearing in the 'cfl3d.out' file:

```
WARNING ... negative volume at i,j,k= 1 514 2 block 1 not stopping! WARNING ... negative volume at i,j,k= 1 515 2 block 1 not stopping!
```

• The majority of negative volumes appear to be in block 1. By plotting the control point output it is clear that elements around the leading edge slat are not well defined, and probably causing poorly defined (singular) macro-elements in that region.

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Example 2: 2D Flap rotation

- The first step in solving this problem is to observe that the file 'meshdef.inp' has been created.
- This file contains the control points that were created by the code.
- Contents of this file can be pasted into the input and customized as needed.
- Since negative volumes occurred in block 1 we will add to the control points in that block.

Contents of 'meshdef.inp':

GRID NIND NJND NKND	
1 2 11 2	

1 2	

1 49 50 137 273 317 473 609 696 697	
745	

1 57	
GRID NIND NJND NKND	
2 2 11 2	

1 2	

1 49 50 137 145 281 325 461 548 549	
597	

1 89	
GRID NIND NJND NKND	
3 2 8 2	

1 2	

1 49 50 121 129 216 217 265	

1 65	
GRID NIND NJND NKND	
4 2 10 2	

1 2 ***********************************	
1 49 50 121 257 413 549 636 637 685	

1 33	

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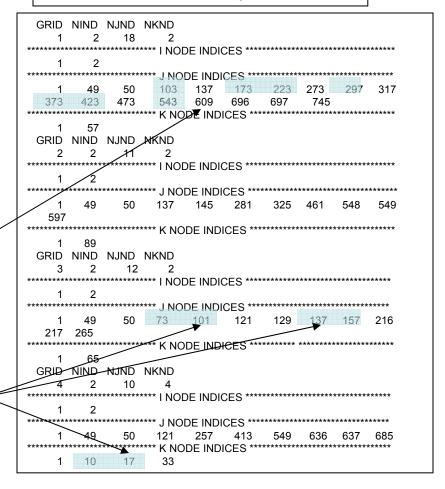
Example 2: 2D Flap rotation

- These additional points have been chosen simply to fill in gaps in the control point distribution.
- This customized input is pasted into the input file, and *nskip* set to 4.

Points added that remove the negative volumes in block 1

Points added to better define the flap region

Contents of 'meshdef.inp' customized:



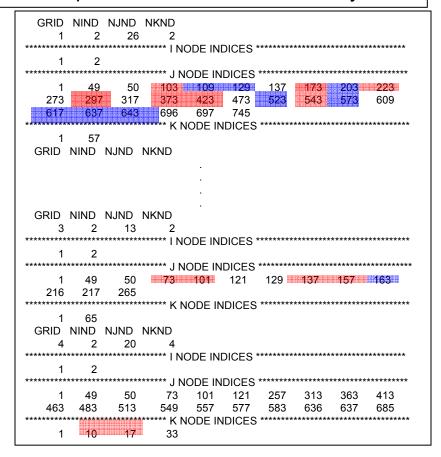


Example 2: 2D Flap rotation

Control point indices the code actually uses:

This is the data output into the new file 'meshdef.inp' after the code is rerun. This file is printed out because new points have been added by the code in addition to points added by the user.

- Control points added by user
- by the code to maintain
 1-1 blocking interface
 continuity

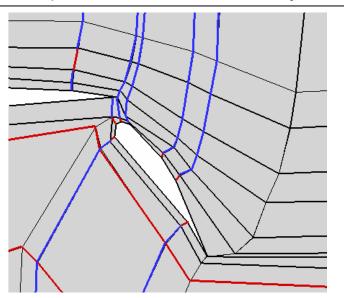


Example 2: 2D Flap rotation



- Control point lines added by the user
- by the code to maintain continuity at 1-1 blocking interfaces

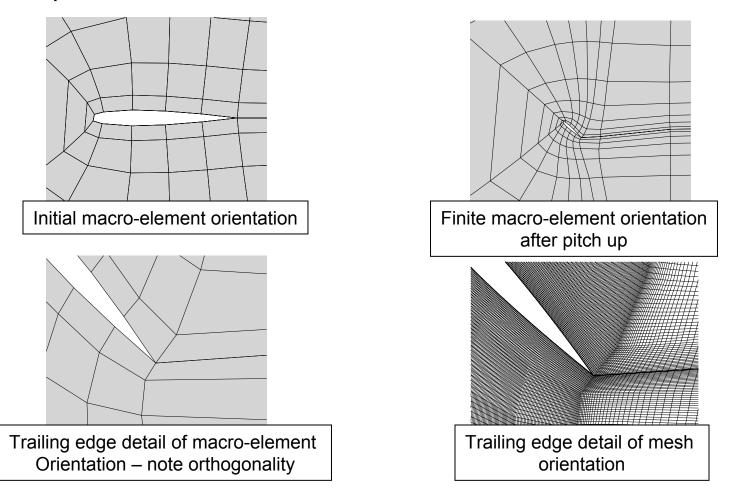
Control point indices the code actually uses:



With these new control points, the code runs robustly with no negative volumes for both the exponential decay and finite macro-element methods for a range of parameter values. Note that the region just aft of the flap retains grid quality better using the finite macro-element method than did the original.



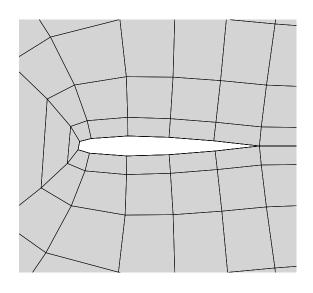
Example 3: 2D airfoil rotation with finite macro-element method



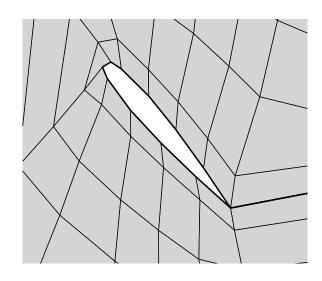
From Bartels, R. E., "Finite Macro-Element Mesh Deformation in a Structured Multi-Block Navier-Stokes Code," NASA/TM-2005-213789, July 2005.

Example 3: 2D airfoil rotation with exponential decay method



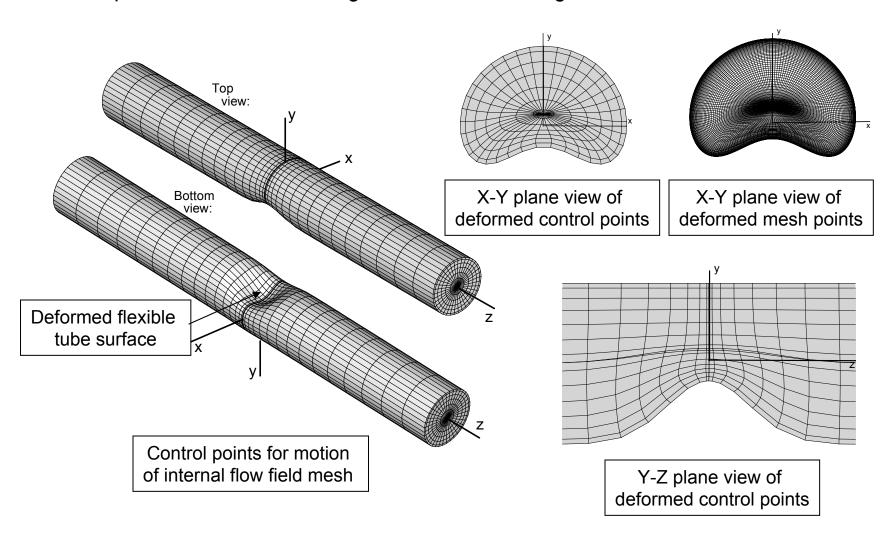


Initial control point orientation



Control point orientation after pitch up, β_2 = 2, α_2 = .005

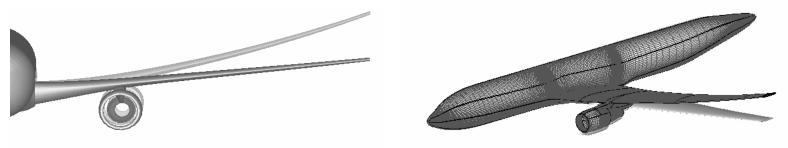
Example 4: Internal flow through a flexible tube using the finite macro-element method

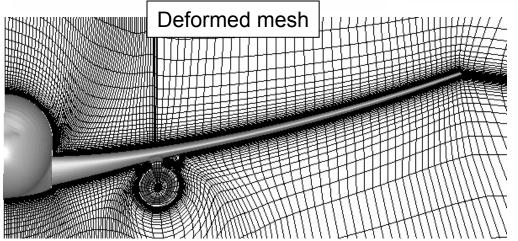




Example 5: Transport wing bending using the Exponential Decay Method

Initial and deformed geometry





This example used mesh movement Option 2 (*isktyp* = -1, *nskip* = 0)

Geometric conservation law

In general the equations computed are

$$\frac{1}{J}\frac{\partial Q}{\partial t} = R(Q)$$

where

 $egin{array}{ll} Q & & - \mbox{ solution vector} \\ J & & - \mbox{ Jacobian of the grid transformation} \end{array}$

R(O) - right hand side composed of spatial flux terms

For steady and unsteady computations:

$$R(Q) = -\left[\frac{\partial (F - F_{v})}{\partial \xi} + \frac{\partial (G - G_{v})}{\partial \eta} + \frac{\partial (H - H_{v})}{\partial \zeta}\right]$$

where

F,G,H - inviscid fluxes F_{v},G_{v},H_{v} - viscous fluxes



Geometric conservation law

For unsteady deforming mesh computations there is an additional term:

$$R(Q) = -\left[\frac{\partial(F - F_v)}{\partial \xi} + \frac{\partial(G - G_v)}{\partial \eta} + \frac{\partial(H - H_v)}{\partial \zeta}\right]$$

$$+ Q\left[\frac{\partial}{\partial t}\left(\frac{1}{J}\right) + \frac{\partial}{\partial \xi}\left(\frac{\xi_t}{J}\right) + \frac{\partial}{\partial \eta}\left(\frac{\eta_t}{J}\right) + \frac{\partial}{\partial \zeta}\left(\frac{\zeta_t}{J}\right)\right]$$
Geometric Conservation Law (GCL), due to grid volume

change

The implication of this is that a computation using rigid grid motion *may* perform somewhat differently than a deforming grid solution with the same time step size, number of sub-iterations and CFL number. However, the two *fully converged* solutions will be the same. See Bartels, R. E., "Mesh and Solution Strategies and the Accurate Computation of Unsteady Spoiler and Aeroelastic Problems," *Journal of Aircraft*, Vol. 37, No. 3, May 2000, pp. 521-529.

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Multiple types of coupled motion

Consider the example of wing plunge combined with control surface rotation. Since the control surface rotation is about a point fixed on the larger moving wing surface, coupling of the two motions will be required. There are two ways to perform this coupled motion:

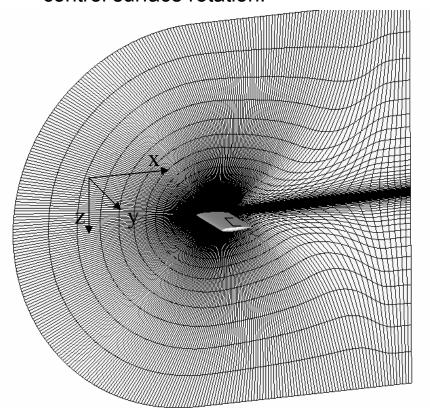
- Coupling control surface rotation and wing translation combined using mesh deformation.
- 2. Coupling control surface rotation using mesh deformation with rigid grid translation.

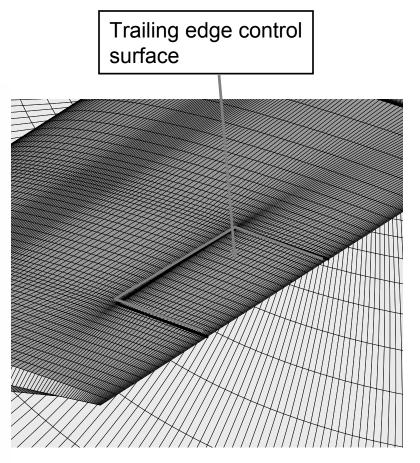
Although these two approaches result in identical wing surface motion, off body grid motion will be much different.



Example: Control surface rotation plus wing plunging

As an example consider the wing shown having both wing plunge plus control surface rotation:







Example: Multi-motion using deforming mesh

The following unsteady input file performs the wing plunging with control surface rotation using deforming mesh:

```
input/output files:
 wbgrid.cfl
 plot3dq.bin
 plot3dq.bin
 cfl3d.out
 cfl3d.res
 cfl3d.turres
 cfl3d.blomax
 cfl3d.out15
 cfl3d.prout
 cfl3d.out20
                                                                                       Note that iunst = 2 since
 ovrlp.bin
                                                                                       deforming mesh is used
 patch.bin
 restart.bin
NASA Langley BACT Model: NACA 0012 af, AR=1.5 wing, .75TE Flap
                                 ReUe
                                         Tinf,dR
                                                    ialph
                                                           ihstry
   Mach
            alpha
                      beta
                                         486.00
 0.82000 0.00000 0.00000 0.236E+07
                                                       1
     sref
              cref
                       bref
                                  xmc
                                           ymc
                                                     zmc
                                        0.00000 0.00000
                              0.25000
   1.000 1.00000 1.00000
                                                  cff tau
      dt
              irest
                     iflagts
                                 fmax
                                           iunst
                                                2.00000
 0.04000
                      3000
                               1.00000
                                                      i2d
    ngrid nplot3d
                     nprint
                                nwrest
                                            ichk
                                                           ntstep
                                                                     ita
                                 1000
                                                                     -2
```



Example: Multi-motion using deforming mesh

	٠,	ivisc(j)	()	iforce	iadvance	•	
	5	5	5	1	0	2 0	
					kdim	, ,	idim
					73		73
		klamhi	klamlo	jlamhi	jlamlo	ilamhi	ilamlo
		.0	. 0	.0	.0		. 0
ke	je	ie	ks	js	is	igridc	inewg
0	0	0	0	0	0	0	0
		iflim(k)		. ,	idiag(k)	idiag(j)	idiag(i)
		3	3	3	1	1	1
			rkap0(j)			fds(j)	ifds(i)
		0.3333		0.3333	1	1	1
iovrlp	nbckdim		nbcjdim	nbcj0	nbcidim	nbci0	grid
0	1	5	1	1	1	1	1
ndata	kend	ksta	jend	jsta	bctype	segment	i0: grid
0	73	1	345	1	1001	1	1
ndata	kend	ksta	jend	jsta	bctype	segment	idim: grid
0	73	1	345		1002	1	1
ndata	kend	ksta	iend	ista	bctype	segment	j0: grid
0	73	1	73	1	1003	1	1
ndata	kend	ksta	end		bctype	segment	jdim: grid
0	73	1	73	1	1003	1	1
ndata	jend	jsta	iend	ista	bctype	segment	k0: grid
0	33	1	49		0	1	1
2	313	33	49	1	2004	2	1
						cq	tw/tinf
						0.00000	0.00000
0	345	313	49	1	0	3	1
0	173	1	73	49	0	4	1
0	345	173	73	. •	0	5	1
ndata	jend	jsta	iend		bctype	segment	kdim: grid
0	345	1	73	1	1003	1	1

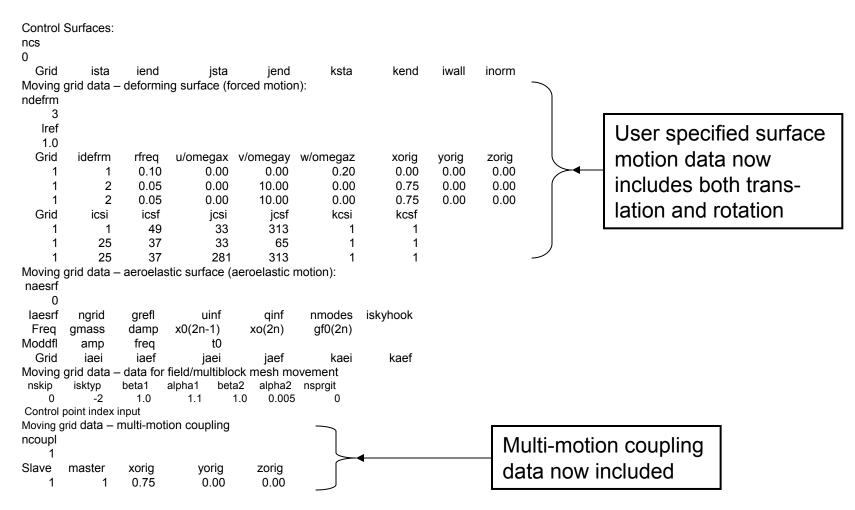


Example: Multi-motion using deforming mesh

mseq	mgflag	iconsf	mtt 0	ngam 2						
inno	2 20000(1)	ا ۱ (۲)ممممر	-	_	nnoor(1)	2000r(2)	opoor(2)			
		epsssc(2)			epsssr(1)					
0	0.3000	0.3000	0.3000	0	0.3000	0.3000	0.3000			
ncyc	mglevg	nemgl	nitfo							
8	3	0	0							
mit1	mit2	mit3	mit4	mit5						
1	1	1								
1-1 block	ing data:									
nbli										
2										
number	grid	ista	jsta	ksta	iend	jend	kend	isva1	isva2	
1	1	1	1	1	49	33	1	1	2	
2	1	49	1	1	73	173	1	1	2	
number	grid	ista	ista	ksta	iend	iend	kend	isva1	isva2	
1	1	1	345	1	49	313	1	1	2	
2	1	49	345	1	73	173	1	1	2	
_	rface date		343		75	173		'		
ninter	erface data	a.								
0										
plot3d ou	•									
grid	iptyp	ista	iend	iinc	jsta	jend	jinc	ksta	kend	kinc
1	0	1	49	1	1	345	1	1	1	1
movie										
0										
print out:										
grid	iptyp	ista	iend	iinc	jsta	jend	jinc	ksta	kend	kinc
1	0	1	49	1	, 1	345	. 1	1	1	1
	U	•	.0		•	310			•	•



Example: Multi-motion using deforming mesh





Example: Multi-motion using deforming mesh

Focusing on the user specified motion input:

Moving grid data – deforming surface (forced motion): ndefrm 3 Iref 1.0 Grid idefrm rfreq u/omegax v/omegay w/omegaz xorig yorig zorig 0.00 0.00 0.10 0.00 0.20 0.00 0.00 The new lines prescribe 0.00 1 0.05 0.00 10.00 0.00 0.75 0.00 the motion of the wing 0.00 2 1 0.05 0.00 10.00 0.00 0.75 0.00 Grid icsi icsf jcsf kcsi kcsf jcsi surface 49 33 313 1 25 37 33 65 1 25 1 37 281 313

Note that idefrm = 1, which corresponds to translational motion.



Example: Multi-motion using deforming plus rigid grid motion

The following unsteady input file performs the wing plunging using rigid grid translation and control surface rotation using deforming mesh:

```
input/output files:
 wbgrid.cfl
 plot3dq.bin
 plot3dq.bin
 cfl3d.out
 cfl3d.res
 cfl3d.turres
 cfl3d.blomax
 cfl3d.out15
 cfl3d.prout
 cfl3d.out20
                                                                                       Note that iunst = 3, for
 ovrlp.bin
                                                                                       deforming mesh plus
 patch.bin
 restart.bin
                                                                                       rigid grid motion
NASA Langley BACT Model: NACA 0012 af, AR=1.5 wing, .75TE Flap
                                ReUe
                                         Tinf,dR
                                                    ialph
                                                           ihstry
   Mach
            alpha
                      beta
                                         486.00
 0.82000 0.00000 0.00000 0.236E+07
                                                      1
     sref
              cref
                      bref
                                  xmc
                                           ymc
                                                    zmc
                              0.25000
                                        0.00000 0.00000
   1.000 1.00000 1.00000
                                                  cff tau
      dt
             irest
                     iflagts
                                 fmax
                                           iunst
                                                2.00000
 0.04000
                      3000
                               1.00000
                                                      i2d
    ngrid nplot3d
                     nprint
                                nwrest
                                            ichk
                                                           ntstep
                                                                    ita
                                 1000
                                                                     -2
```



Example: Multi-motion using deforming plus rigid grid motion

• • • • • • • • • • • • • • • • • • • •	c(k)
2 0 0 1 5 5	5
idim jdim kdim	
73 345 73	
ilamlo ilamhi jlamlo jlamhi klamlo klamhi	
0 0 0 0 0	
inewg igridc is js ks ie	je ke
0 0 0 0 0	0 0
idiag(i) idiag(j) idiag(k) iflim(i) iflim(j) iflim(k)	
1 1 1 3 3 3	
ifds(i) fds(j) ifds(k) rkap0(i) rkap0(j) rkap0(k)	
1 1 1 0.3333 0.3333 0.3333	
grid nbci0 nbcidim nbcj0 nbcjdim nbck0 nbck	dim iovrlp
1 1 1 1 5	1 0
i0: grid segment bctype jsta jend ksta k	end ndata
1 1 1001 1 345 1	73 0
idim: grid segment bctype jsta jend ksta ko	end ndata
1 1 1002 1 345 1	73 0
j0: grid segment bctype ista iend ksta k	end ndata
1 1 1003 1 73 1	73 0
jdim: grid segment bctype ista end ksta k	end ndata
1 1 1003 1 73 1	73 0
k0: grid segment bctype ista iend jsta je	end ndata
1 1 0 1 49 1	33 0
1 2 2004 1 49 33	313 2
tw/tinf cq	
0.00000 0.00000	
1 3 0 1 49 313	345 0
1 4 0 49 73 1	173 0
1 5 0 49 73 173	345 0
kdim: grid segment bctype ista iend jsta je	end ndata
	345 0

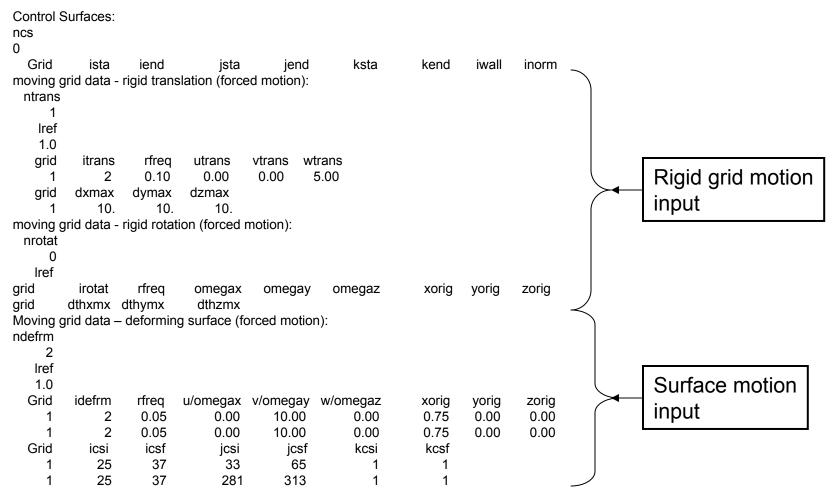


Example: Multi-motion using deforming plus rigid grid motion

0 ncyc 8	0.3000 mglevg 3	iconsf 1 epsssc(2) 0.3000 nemgl 0	0.3000 nitfo 0	0	0.3000	epsssr(2) 0.3000	epsssr(3) 0.3000			
mit1	mit2	mit3	mit4	mit5	•••					
1 1-1 block nbli 2	1 ing data:	1								
number	grid	ista	jsta	ksta	iend	jend	kend	isva1	isva2	
1	1	1	1	1	49	33	1	1	2	
2	1	49	1	1	73	173	1	1	2	
number	grid	ista	jsta	ksta	iend	jend	kend	isva1	isva2	
1	1	1	345	1	49	313	1	1	2	
2	1	49	345	1	73	173	1	1	2	
patch inter ninter 0	erface data	a:								
plot3d ou	tput:									
grid	iptyp	ista	iend	iinc	jsta	jend	jinc	ksta	kend	kinc
1	0	1	49	1	1	345	1	1	1	1
movie 0										
print out:		:-4-	: d		:-4-			14	1	1.5
grid 1	iptyp 0	ista 1	iend 49	iinc 1	jsta 1	jend 345	jinc 1	ksta 1	kend 1	kinc 1

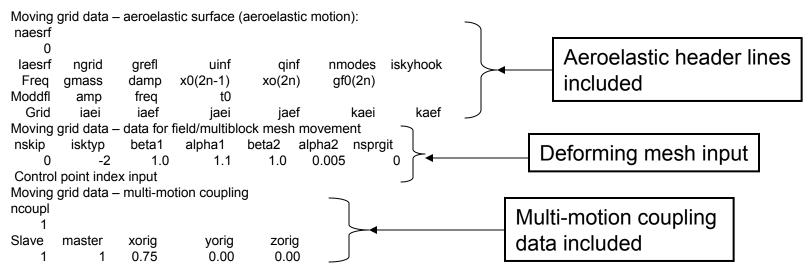


Example: Multi-motion using deforming plus rigid grid motion





Example: Multi-motion using deforming plus rigid grid motion



Note: CFL3D does not allow initiating new kinds of motion upon restarts. Therefore if an initial deforming mesh computation is performed to reach an equilibrium before initiating a combined rigid and moving (deforming) control surface computation, the option iunst = 3 must be used from the start (that is after an initial steady state computation with dt < 0), with control surface motion set to zero.

Overview



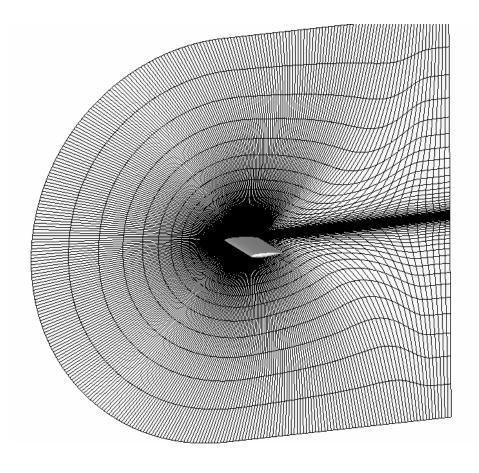
- CFL3D has the capability to perform both static and dynamic aeroelastic analysis. In this analysis the fluid and structure interact through a time marching simulation (e.g. flutter analysis, etc...)
- All aeroelastic and modal analyses are performed by running the code in unsteady mode
- CFL3D performs aeroelastic analysis for a linear structure modally and a linear and nonlinear structure when loosely coupled with either NASTRAN or ABAQUS finite element analyses.
- For modal analysis, the equations of structural dynamics must be decoupled modally
 - Eigenvalue analysis is required prior to running CFD to obtain frequencies, generalized masses and mode shapes.
 - A preprocessing step projecting the mode shapes onto the CFD surface grids is required.
 - The code reads the modal data projected onto the CFD surfaces in the file 'aesurf.dat'. This
 file must be contained in the directory in which the executable resides.
- CFL3D also has the capability to perform unsteady deforming body analysis using mode shapes. In this mode the user specifies modal motion (e.g. control surface rotation, wing plunge oscillation, etc...) in the aeroelastic input section

Example of an aeroelastic model



Consider the Benchmark Active
Controls Technology (BACT)
aeroelastic model shown. The
model has pitch and plunge
aeroelastic degrees of freedom. The
model parameters are:

$$M_T = 5.966$$
 slugs
 $S_{\alpha} = 0.01420$ slug-ft
 $I_{\alpha} = 2.8017$ slug-ft2
 $K_h = 2659$ lb/ft
 $K_a = 2897$ lb-ft/rad



NA

Example of an aeroelastic model

The coupled equations of structural dynamics are

$$\begin{bmatrix} M_T & S_{\alpha} \\ S_{\alpha} & I_{\alpha} \end{bmatrix} \left\{ \ddot{\zeta} \right\} + \begin{bmatrix} K_h & 0 \\ 0 & K_{\alpha} \end{bmatrix} \left\{ \zeta \right\} = q_{\infty} \left\{ \iint c_p(x^*, y^*) dx^* dy^* \\ \iint c_p(x^*, y^*) (x^*_{ea} - x^*) dx^* dy^* \right\}$$

where ζ_1 is plunge (h) and ζ_2 is pitch (α). Eigen-analysis of this system yields the frequencies

$$\omega_h = 21.1113283 \ rad \ / \sec (3.36 \ Hz)$$

$$\omega_{\alpha} = 32.1564455 \ rad \ / \sec (5.12 \ Hz)$$

Example of an aeroelastic model



Using the eigenvectors

$$\phi = \begin{bmatrix} \varphi_{11} & \varphi_{12} \\ \varphi_{21} & \varphi_{22} \end{bmatrix} = \begin{bmatrix} 0.409404775 & 0.0024991919 \\ 0.001571926 & -0.5974345042 \end{bmatrix}$$

the generalized masses are obtained

$$m_1 = 1.0000000000$$

$$m_2 = 1.0000000000$$

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Example of an aeroelastic model

... and the decoupled equations of structural dynamics

$$\begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \{ \ddot{q} \} + \begin{bmatrix} \omega_{1}^{2} & 0 \\ 0 & \omega_{2}^{2} \end{bmatrix} \{ q \} = \mathbf{M}^{-1} \phi^{\mathsf{T}} \mathbf{q}_{\infty} \left\{ \iint c_{p}(x^{*}, y^{*}) dx^{*} dy^{*} \\ \iint c_{p}(x^{*}, y^{*}) (x^{*}_{ea} - x^{*}) dx^{*} dy^{*} \right\}$$

where

$$\zeta = \phi \, q \quad M = \begin{bmatrix} m_1 & 0 \\ 0 & m_2 \end{bmatrix} \quad \begin{bmatrix} m_1 \omega_1^2 & 0 \\ 0 & m_2 \omega_2^2 \end{bmatrix} = \phi^T \begin{bmatrix} K_h & 0 \\ 0 & K_\alpha \end{bmatrix} \phi$$

Carrying through the multiplication on the right-hand side, we have

$$\begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \{ \ddot{q} \} + \begin{bmatrix} \omega_{1}^{2} & 0 \\ 0 & \omega_{2}^{2} \end{bmatrix} \{ q \} = \mathbf{M}^{-1} \mathbf{q}_{\infty} \left\{ \iint_{\mathbf{p}} c_{p}(x^{*}, y^{*}) \{ \boldsymbol{\varphi}_{11} + \boldsymbol{\varphi}_{21}(x^{*}_{ea} - x^{*}) \} dx^{*} dy^{*} \right\}$$

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Example of an aeroelastic model

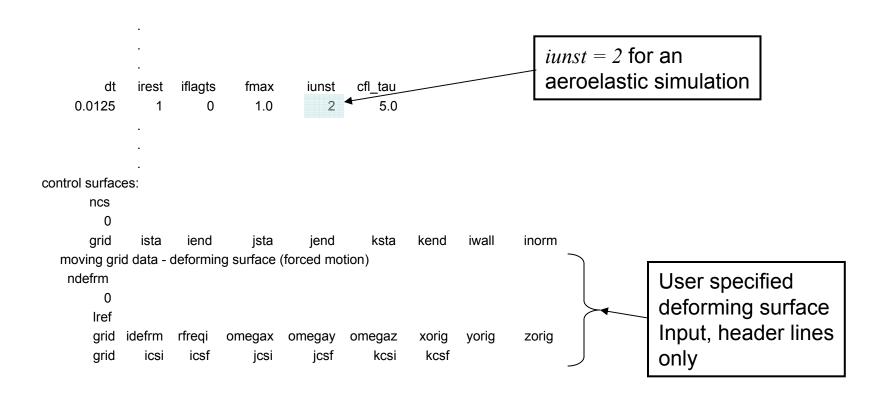
The mode shapes that are input into CFL3D are revealed by the last equations

$$\begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \{ \ddot{q} \} + \begin{bmatrix} \omega_{\rm l}^2 & 0 \\ 0 & \omega_{\rm 2}^2 \end{bmatrix} \{ q \} = \mathbf{M}^{\text{-1}} \mathbf{q}_{\infty} \begin{cases} \iint c_p(x^*, y^*) \{ \varphi_{11} + \varphi_{21}(x^*_{\text{ea}} - x^*) \} dx^* dy^* \\ \iint c_p(x^*, y^*) \{ \varphi_{12} + \varphi_{22}(x^*_{\text{ea}} - x^*) \} dx^* dy^* \end{cases}$$
 First mode shape, $\Phi_{\rm z,1}$

These can be used to create the modal shape projected to each wing surface grid point for input into CFL3D. Note that x^* and y^* are in the same units as the structural model.

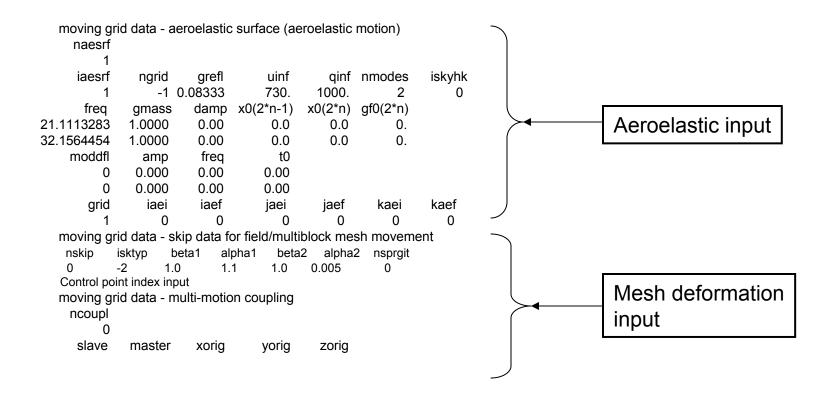
Aeroelastic input





Aeroelastic input

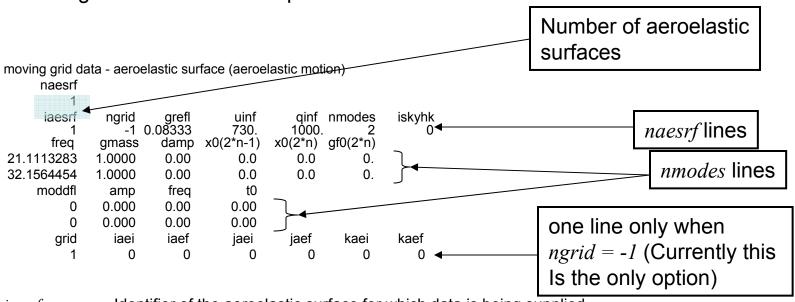




Aeroelastic input



Focusing on the aeroelastic input section:



iaesrf

- Identifier of the aeroelastic surface for which data is being supplied

ngrid

- Number of surface segments that make up this aeroelastic surface

nmodes

- Number of modes to be modeled in CFL3D

iskyhk

- Not currently used, any value will serve as a placeholder

uinf

- Free-stream velocity, in the same units as the equations of structural dynamics

qinf

- Dynamic pressure, in the same units as the equations of structural dynamics

grefl

- Conversion from CFD grid units to structural equation units.

Aeroelastic input



Regarding the input parameter *grefl*, consider the equations of structural dynamics for the pitch/plunge example:

$$\begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \{ \ddot{q} \} + \begin{bmatrix} \omega_1^2 & 0 \\ 0 & \omega_2^2 \end{bmatrix} \{ q \} =$$

$$\begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \{ \ddot{q} \} + \begin{bmatrix} \omega_1^2 & 0 \\ 0 & \omega_2^2 \end{bmatrix} \{ q \} = M^{-1} q_{\infty} \left\{ \iint_{c_p(x^*, y^*)} \Phi_{z,1} dx^* dy^* \right\}$$

The actual equations solved in CFL3D are:

Lengths in structural model units

$$\begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \left\{ \ddot{q} \right\} + \begin{bmatrix} \omega_1^2 & 0 \\ 0 & \omega_2^2 \end{bmatrix} \left\{ q \right\} = \underbrace{\operatorname{grefl}^2}_{} \operatorname{M^{-1}} \operatorname{q}_{\infty} \left\{ \iint_{} c_p(x,y) \Phi_{z,1} \, dx \, dy \right\} \\ \iint_{} c_p(x,y) \Phi_{z,1} \, dx \, dy \right\}$$
 By definition:
$$\operatorname{grefl} = \sqrt{S_{AE} \, / \, S_{CFD}}$$
 Lengths in CFD grid units

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Aeroelastic input

In the present example the structural equations are in units of feet, while the CFD grid is in units of inches. **Note that the aspect ratios of the original and rescaled model must be identical**. Conversion for the present example can be obtained from

$$grefl = \sqrt{S_{AE}/S_{CFD}} = \sqrt{\frac{1}{144}} \approx 0.08333 ft/grid unit$$

This is the *grefl* parameter that would be entered in the aeroelastic input section.

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Modal form of the equations

Consider the decoupled equations of structural dynamics for N (or nmodes in the input) modes

$$\begin{bmatrix} 1 & 0 & 0 \\ 0 & \ddots & 0 \\ 0 & 0 & 1 \end{bmatrix} \{ \ddot{q} \} + \begin{bmatrix} 2\omega_{1}\zeta_{1} & 0 & 0 \\ 0 & \ddots & 0 \\ 0 & 0 & 2\omega_{N}\zeta_{N} \end{bmatrix} \{ \dot{q} \} + \begin{bmatrix} \omega_{1}^{2} & 0 & 0 \\ 0 & \ddots & 0 \\ 0 & 0 & \omega_{N}^{2} \end{bmatrix} \{ q \}$$

$$= \begin{bmatrix} m_{1}^{-1} & 0 & 0 \\ 0 & \ddots & 0 \\ 0 & 0 & m_{N}^{-1} \end{bmatrix} \{ Q \}$$

where q is the modal variable vector and Q is the generalized force vector, each of length N. ω_1 ,..., ω_N are the natural frequencies of each structural mode in radians, and m_1 ,..., m_N are the generalized masses.

Modal form of the equations



CFL3D input definitions as they relate to the modal equations of structural dynamics are as follows:

$$gmass(1) = m_1, \quad \cdots, \quad gmass(N) = m_N$$

$$freq(1) = \omega_1, \quad \cdots, \quad freq(N) = \omega_N$$

$$damping(1) = \zeta_1, \quad \cdots, \quad damping(N) = \zeta_N$$

$$x0(1) = q_{1 init}, \cdots, \quad x0(2*N-1) = q_{N init}$$

$$x0(2) = \dot{q}_{1 init}, \cdots, \quad x0(2*N) = \dot{q}_{N init}$$

$$gf(0) = Q_{1 init}, \cdots, \quad gf(0) = Q_{N init}$$

Units for frequency is radians/time (usually time scale is seconds for the structural dynamics equations).

Aeroelastic input



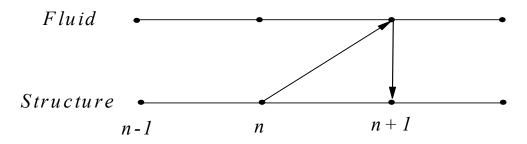
- $x\theta(2*n-1)$ is the initial generalized displacement of the mode; will override the value in the restart file (if restarting) when $x\theta(2*n-1)$ is nonzero. Otherwise, it will not override the restart value. This allows the mode to be perturbed for excitation of aeroelastic dynamic response after a static aeroelastic starting solution has been performed.
- $x\theta(2*n)$ is the initial generalized velocity of the mode; will override the value in the restart file (if restarting) when $x\theta(2*n)$ is nonzero. Otherwise, it will not override the restart value. This allows the mode to be perturbed for excitation of aeroelastic dynamic response after a static aeroelastic starting solution has been performed.
- gf0(2*n) is the generalized force offset to include for the mode. This value is included in CFL3D computation of generalized force in the following way for mode n = 1 to nmodes:

$$Q_n = q_{\infty} \ grefl^2 \left\{ \iint c_p \ \vec{\Phi}_n \cdot d\vec{s} \right\} - gf \ 0 (2*n)$$
Value from input



Method of integrating the fluid/structure coupling based on the state transition matrix solution

The structural dynamic equations are written in state space form and solved using the state transition matrix solution and a predictor/corrector scheme. The fluid/structure cycling is shown below:



Predictor step:

$$\widetilde{x}^{n+1} = \Theta x^n + \frac{1}{2} \Theta_i \left(3 \overline{Q}^n - \overline{Q}^{n-1} \right)$$

After the predictor step, the structure surface is moved using the predicted modal solution \tilde{x}^{n+1} , the flow field is converged to the solution at time step n+1, and the new generalized force \overline{Q}^{n+1} is computed.

Corrector step:
$$x^{n+1} = \Theta x^n + \frac{1}{2} \Theta_i \left(\overline{Q}^{n+1} + \overline{Q}^n \right)$$



Method of integration and fluid/structure coupling

The parameter Θ is the state transition matrix. The parameter Θ_i is the discrete integration of the state input, and \overline{Q}^n is the state vector containing the generalized aerodynamic force at time step n. The solution and generalized

force state vectors are

$$x = egin{cases} q_1 \ \dot{q}_1 \ \vdots \ q_N \ \dot{q}_N \end{pmatrix} \qquad , \qquad \overline{Q} = egin{cases} 0 \ Q_1 \ \vdots \ 0 \ Q_N \end{pmatrix}$$

The value \widetilde{x} is the intermediate state solution used to update the mesh.

This method uses a second order backward differencing of the fluid/structure coupling. Combined with second order backward differenced fluid dynamics solver and the second order backward differencing of the mesh time metrics, the overall scheme is second order accurate. For the original method see: Edwards, J. W., Bennett, R. M., Whitlow Jr., W., Seidel, D. A., "Time-Marching Transonic Flutter Solutions Including Angle-of-Attack Effects," *Journal of Aircraft*, 20 (1983), pg. 899-906.

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Modal surface input

- Currently CFL3D assumes that the aeroelastic surface comprises all boundary segments with the boundary condition types 1005, 1006, 2004, 2014 or 2016.
- Note that the boundary condition 1001 is not considered an aeroelastic surface. Therefore, if a symmetry plane is required to deform with a pitching wing, it must be treated as an inviscid wall boundary (1005 or 1006)
- The modal input file *aesurf.dat* must have modal data for a given surface point in free field ascii format (no commas) with $\Phi_{x,n}$, $\Phi_{y,n}$, $\Phi_{z,n}$ modal deflections at each surface point for each mode n.

Format of the modal surface input

The following ordering is required:

Segment limits defined in boundary condition input

$$j=1$$
 surface:
 $\Phi_{x,n}(i,j,k)$ $\Phi_{y,n}(i,j,k)$ $\Phi_{z,n}(i,j,k)$

$$k = ksta$$
 to $kend$, $i = ista$ to $iend$, repeat $nseg$ times

$$j = jdim$$
 surface:

$$\Phi_{x,n}(i,j,k)$$
 $\Phi_{v,n}(i,j,k)$ $\Phi_{z,n}(i,j,k)$

$$k = ksta$$
 to $kend$, $i = ista$ to $iend$, repeat $nseg$ times

$$k = 1$$
 surface:

$$\Phi_{v,n}(i,j,k)$$
 $\Phi_{v,n}(i,j,k)$ $\Phi_{z,n}(i,j,k)$

,
$$j = ista$$
 to $iend$, $i = ista$ to $iend$, repeat $nseg$ times

k = kdim surface:

$$\Phi_{x,n}(i,j,k)$$
 $\Phi_{y,n}(i,j,k)$ $\Phi_{z,n}(i,j,k)$

,
$$j = jsta$$
 to $jend$, $i = ista$ to $iend$, repeat $nseg$ times

i = 1 surface:

$$\Phi_{x,n}(i,j,k)$$
 $\Phi_{y,n}(i,j,k)$ $\Phi_{z,n}(i,j,k)$

, j = jsta to jend , k = ksta to kend , repeat nseg times

i = idim surface:

$$\Phi_{x,n}(i,j,k)$$
 $\Phi_{y,n}(i,j,k)$ $\Phi_{z,n}(i,j,k)$

, j = jsta to jend, k = ksta to kend , repeat nseg times,

Repeat all of the above input for n = 1 to nmodes, repeat ngrid times, repeat naesrf times.

Format of the modal surface input



- The ordering of the aeroelastic surface points must correspond to the order of the points in the CFD grid file read by CFL3D.
- Aeroelastic segments must be input in the same block order as the grid file, and segments must be input in order of ascending indices.
- When creating a multi zonal grid using the utility 'splitter', be aware that the final ordering will generally *not* correspond to the ordering of the unsplit grid. Ordering of the split grid zones can be found in the 'splitter.out' file, from which can be found the required order of the surface grid points for the 'aesurf.dat' file.

Example: Consider a block face that has dimensions kdim = 49, idim = 49 with several aeroelastic segments. If segment 1 has indices k = 33 to 49, i = 13 to 33, and segment 2 has indices k = 1 to 33, then segment 2 must be input first.

Aeroelastic output

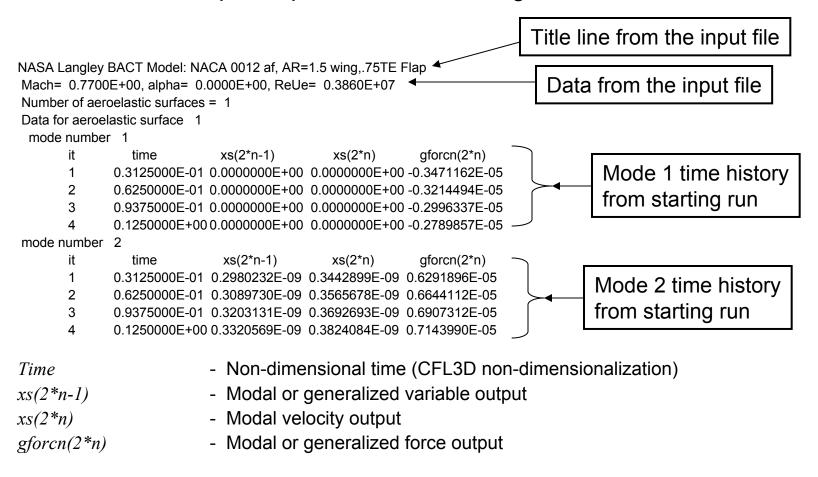


- Aeroelastic time history output is in the file 'genforce.dat'.
- This file is generated if iunst = 2 and aeroelastic surfaces are defined in the input file $(naesrf \neq 0)$.
- After header information, modal response data for each mode is written sequentially.
- Unlike output data in the 'cfl3d.subit_res' file, a complete time history of this data for the entire simulation is retained and written/read to/from restart files and subsequently output to the 'genforce.dat' file.

Aeroelastic output



Consider the example output contained in the 'genforce.dat' file:



Strategy for aeroelastic computations

The following strategies may be used for performing static or dynamic aeroelastic simulations

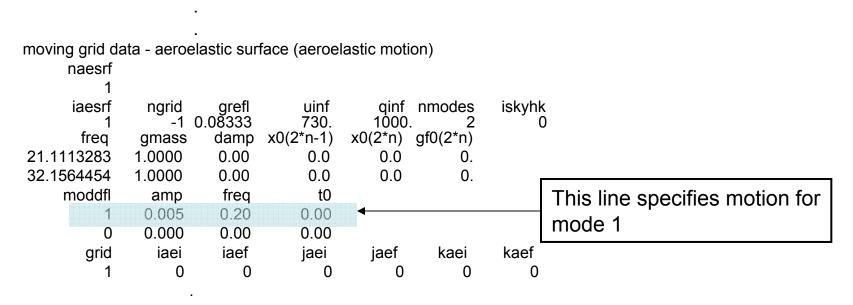
- Static aeroelastic computations can be performed by:
 - Start either from scratch (irest = 0), or restart, after a steady state computation (in which dt < 0, iunst = 0). Starting from scratch is not recommended.
 - Set iunst = 2, dt > 0 and damp = .999999... and perform the computation in a time marching manner to convergence.
- Flutter onset computations can be performed by:
 - Converging a static solution as outlined above.
 - Setting damp to the correct value for the elastic system being modeled.
 - Setting an initial perturbation $x\theta(2*n)$ or $x\theta(2*n-1)$ in the desired mode.*

^{*} If a restart in the middle of a flutter computation is performed, the initial perturbation values from the previous run must be reset to zero at the restart of the new run.

User specified modal motion



The user may specify modal motion within the aeroelastic input (e.g. control surface rotation, wing plunge oscillation, impulse for frequency response, etc...) The following modifications to the aeroelastic input specifies modal motion:



User specified modal motion



moddfl

type of time-varying modal perturbation desired:

- < 0, mode displacement and velocity set to zero
- = 0, no perturbation (solution via the dynamic modal equations)
- = 1, harmonic (sinusoidal) perturbation
- = 2, Gaussian pulse
- = 3, step pulse

A (amp)

amplitude of modal perturbation.

ω_r (freq)

reduced frequency of modal perturbation if moddfl = 1 half-width of Gaussian pulse if moddfl = 2 use any value as a placeholder for moddfl = 0

 t_0 (t0)

time about which Gaussian pulse is centered if moddfl = 2 time of the step pulse if moddfl = 3 use any value as a placeholder for moddfl = 0

User specified modal motion



For harmonic perturbation the modal displacement and velocities for mode *n* are computed in the following way:

$$q_n = A\sin(\omega_r t^*)$$
 , $\dot{q}_n = Ak_r\cos(\omega_r t^*)$

where A = amp, $\omega_r = freq$ in radians per dimensional time, and t^* is dimensional time,

$$t^* = t \operatorname{grefl} / a_{\infty}$$
 , $a_{\infty} = U_{\infty} / M_{\infty}$

 U_{∞} (uinf) is in the aeroelastic input section and M_{∞} is from the main aerodynamic input section. t is CFL3D non-dimensional time.

For a Gaussian pulse the displacement and velocity for mode n are computed with

$$q_n = Ae^{-C[t^*-t_0]^2}, \dot{q}_n = -2CAe^{-C[t^*-t_0]^2}$$

where $C = \log(2)/\omega_r^2$

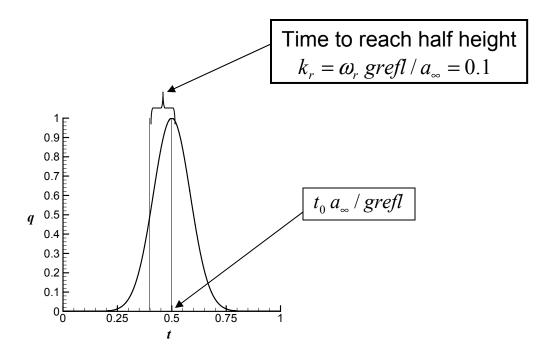
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User specified modal motion

For step pulse the modal displacement and velocities for mode *n* are computed in the following way:



Example: Gaussian modal pulse and time step sizing



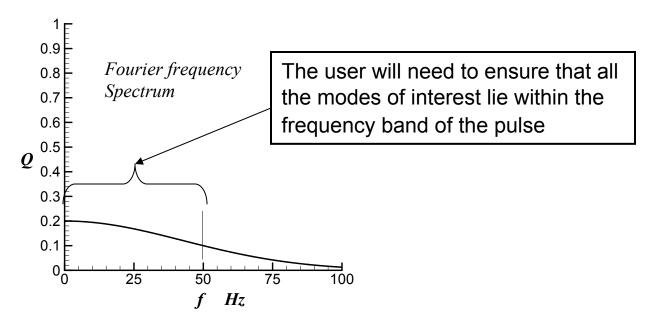
For this example:

$$A=1.0$$
 , $k_r=0.1$, $t_0=0.5 \, grefl/a_{\infty}$
 $C=\log(2)/\omega_r^2$ $t^*=t \, grefl/a_{\infty}$, $a_{\infty}=U_{\infty}/M_{\infty}$

Recommend sizing time step so that there are an absolute *minimum* of 25 time steps within the half life of the pulse ($\Delta t = k_r/25$). In this case we would have $\Delta t = 0.004$.



Example: Shaping and sizing the Gaussian modal pulse

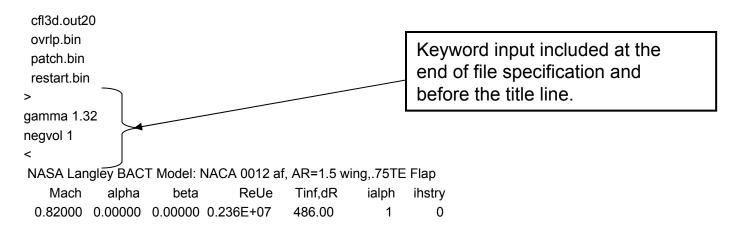


- For a linear response, we will usually want the amplitude as small as possible while staying significantly (say several orders of magnitude) above numerical round off errors.
- Low frequency responses will be very sensitive to the steady convergence of a solution. Therefore, great care must be exercised in adequately converging the steady state if an FRF is the desired outcome.
- The solution is very sensitive to sub-iterative convergence at each time step. A strategy of multiple restarts with different numbers of sub-iterations through the pulse region can reduce the overall run time.

Overview



- There is additional input in CFL3D version 6 that does not fit into an input format consistent with earlier versions of the code. These input parameters have been included as keyword input.
- Keyword input is an optional input specified by lines started by a line with '>'
 and ended with a line containing '<'.
- The following example illustrates how keyword input is included:



Valid Keywords



Physical Properties

Name	Description	Default Value
cbar	Ref. temp. for Sutherland Law	198.6
gamma	Ratio of specific heats	1.4
pr	Prandtl number	0.72
prt	Turbulent Prandtl number	0.90

Limiters

Name	Description	Default Value
atol	Tolerance for detecting singular lines	10 ⁻⁷
epsa_r	Eigenvalue limiter (entropy fix for high Mach flows)	0.0

Valid Keywords



Preconditioning

Name	Description	Default Value
avn	Factor multiplying uref for preconditioning	1.0
cprec	Relative amount of preconditioning	0.0
uref	Limiting velocity for preconditioning	xmach

Specified CL

Name	Description	Default Value
cltarg	Target Cl	99999.
dalim	Limit of alpha change (deg) per update	0.2
icycupdt	Number of cycles between alpha updates (if > 0; if < 0, alpha is never updated)	1
rlxalph	Relaxation factor used to update angle of attack	1.0

Valid Keywords



Name	Description	Default Value
cflturb	Cfl no. for turbl eqns. = cflturb x abs(dt) If cflturb > 0	0
		(model dependent default)
edvislim	Limiter for eddy viscosity in 2-equation	100000.
	turb models; eddy viscosity limited to edvislim times the laminar viscosity	
ibeta8kzeta	flag (0/1) to set beta8 term when using k-enstrophy turbulence model (ivisc=15); 0 = use beta8=0.0 (helps avoid numerical problems); 1 = use beta8=2.3 (available <i>after</i> V6.3)	0
ides	flag (0/1) to perform DES with turbulence model (1) or not (0)	0
cdes	constant associated with DES	0.65
ieasmcc2d	flag (0/1) to turn on 2-D curvature correction when using EASM models (ivisc=8,9,11,12,13,14) (1) or not (0) (available after V6.3)	0
isarc2d	flag (0/1) to turn on 2-D curvature correction when using SA model (ivisc=5) (1) or not (0) (available after V6.3)	0

Valid Keywords



Name	Description	Default Value
sarccr3	value of cr3 parameter in SARC model (available <i>after</i> V6.3)	0.6
ikoprod	flag: 0=use approximate (vorticity-based) turb production term (-2*mut*WijWji) for turb models 6, 7, 10, or 15; 1=use strain-rate based term (2*mut*SijSij); 2=use full production term (ivisc=15 only) (available after V6.3)	0 (vorticity-based production)
isstdenom	flag (0/1): 0=use vorticity term in denominator of eddy viscosity in SST model (#7); 1=use strain term (available after V6.3)	0 (vorticity term)
itaturb	flag (0/1) to control time accuracy of turb. model; 0 for 1st order in time regardless of parameter "ita" for the mean flow; 1 for same order as set by ita	1 (turb. Time accuracy same as mean flow, set via ita)
iturbord	flag controls whether turbulence model advection terms are 1st or 2nd order upwind on RHS (1=1st, 2=2nd) (note: LHS uses 1st order in both cases) (available after V6.3)	1 (1 st order)

Valid Keywords



Name	Description	Default Value
iturbprod	flag: 0=use strain-rate based turb production term (2*mut*SijSij) for EASM turb models 8, 9, 13, or 14; 1=use full production term	0 (strain-rate based term)
nfreeze	Freeze turb. model for nfreeze cycles	0 (not frozen)
nsubturb	Number of iterations of turb model per cycle	1
pklimterm	factor used to limit production of k in 2-eqn turb models (chooses min of Pk and pklimterm*Dk); make this term large for no limiting (available after V6.3)	20.0
tur10 & tur20	turbulent quantity freestream levels < 0 use default value (different for each turb model, see manual Appendix H) =0 use this number as the specified user input value	-1
tur1cut	value that nondimensional epsilon (or omega or enstrophy) is reset to when it tries to drop equal to or below tur1cutlev; if <=0 then no update occurs when value tries to drop equal to or below tur1cutlev (available after V6.3)	1.e-20 for all models except -1 for ivisc=15

Valid Keywords



Name	Description	Default Value
tur2cut	value that nondimensional k is reset to when it tries to drop equal to or below tur2cutlev; if <=0 then no update occurs when value tries to drop equal to or below tur2cutlev (available after V6.3)	1.e-20
tur1cutlev & tur2cutlev	lower levels of nondimensional epsilon (or omega or enstrophy) and k which, when reached, cause the turb quantities to be reset to tur1cut or tur2cut (available <i>after</i> V6.3)	0

Keyword Input Valid Keywords



Deformation/grid motion

Name	Description	Default Value
idef_ss	flag (0/1) to deform volume grid to surface in file	0
	newsurf.p3d	(don't deform)
meshdef	flag (0/1) to bypass flow solution while still computing grid operations such as metrics and volumes; 0 = normal operation; 1 = bypass flow solution (available after V6.3)	0
negvol	flag (0/1) to enable/disable stop if neg. volumes/bad metrics are detected	0 (stop for negative volumes)

Input/output control

Name	Description	Default Value
ibin	flag (0/1) for formatted/unformatted output plot3d files	1 (unformatted)
iblnk	flag (0/1) for un-iblanked/iblanked output plot3d files	1 (iblanked)

Valid Keywords



Input/output control

Name	Description	Default Value
iblnkfr	flag (0/1) for un-iblanked/iblanked fringe points in	1
	plot3d files (overset grids only)	(iblanked)
icgns	flag (0/1) to not use/use CGNS files*	0 (don't use CGNS files)
ip3dgrad	flag (0/1) for solution/derivative data output to	0
	plot3d q file (complex code only)	(solution to q file)
irghost	flag to read ghost-cell data from restart file (1) or	1
	not (0); V5 restart files and Beta V6 restart files do not contain ghost-cell data; newer V6 restart files	(read ghost-cell data)
	do	
iwghost	flag to write ghost-cell data to restart file (1) or not	1
	(0); V5 restart files and Beta V6 restart files do not contain ghost-cell data; newer V6 restart files do	(write ghost-cell data)

Valid Keywords



Input/output control

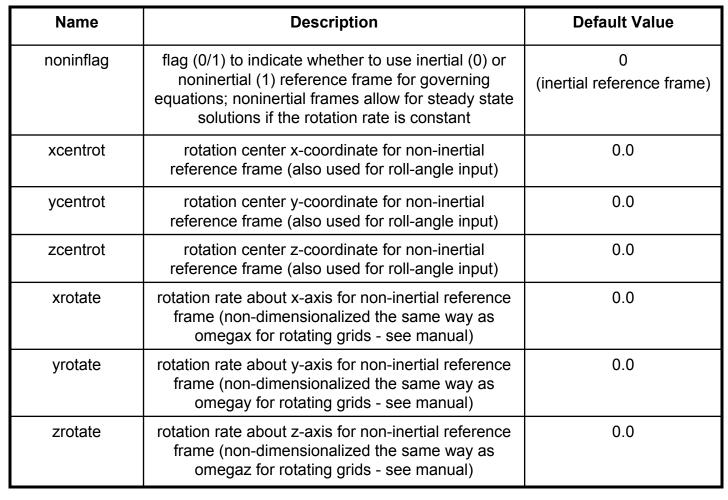
Name	Description	Default Value
itime2read	flag (0/1) to skip/read 2nd order (in time) turbulence terms and dt in restart file: need to skip if using an older time-accurate-with-2nd-order-time restart file	1 (read 2 nd order time turbulence terms and dt)
iteravg	flag to store iteration-averaged conserved variables in PLOT3D files: 0 = no averaging or storage 1 = start averaging now 2 = continue averaging from previous run	0

Memory management

Name	Description	Default Value
memadd	memadd additional memory (in words) added to work array (in case sizer underestimates)	0
		(no addition to work)
memaddi	additional memory (in words) added to iwork array (in case sizer underestimates)	0
		(no addition to iwork)

Valid Keywords

Reference frame





Valid Keywords



Reference frame

Name	Description	Default Value
xrotrate_img	complex perturbation to rotation rate about x-axis for non-inertial reference frame, for computing rate derivatives	0.0
yrotrate_img	complex perturbation to rotation rate about y-axis for non-inertial reference frame, for computing rate derivatives	0.0
zrotrate_img	complex perturbation to rotation rate about z-axis for non-inertial reference frame, for computing rate derivatives	0.0

Other

Name	Description	Default Value
alpha_img	Imaginary perturbation to alpha	0.0
beta_img	Imaginary perturbation to beta	0.0
geom_img	Imaginary perturbation to grid	0.0

Valid Keywords



Other

Name	Description	Default Value
reue_img	Imaginary perturbation to unit Re	0.0
surf_img	Imaginary perturbation to surface grid	0.0
tinf_img	Imaginary perturbation to Tinf	0.0
xmach_img	Imaginary perturbation to Mach no.	0.0
iaxi2plane	flag for use with particular axisymmetric cases (for which i2d=0 and idim=2); if iaxi2plane = 1, the time step based on CFL number is modified so it does not depend on the i-direction metrics (available after V6.3)	0 (no mods to time step)
ifullns	flag (0/1) to specify inclusion of cross-derivative terms; 0 = thin-layer N-S; 1 = full N-S (available after V6.3)	0
ivolint	flag (0/1) to use approximate/exact one-to-one boundary volumes (0 emulates V5.0)	1 (exact volumes)
roll_angle	x-axis roll angle (deg) "+" is clockwise viewed from "- x" (left roll to pilot) (grid is rotated to this angle)	0.0

Overview



- Message Passing Interface (MPI) protocol is used for parallelization of CFL3D
- MPI parallelizes by parceling out grid blocks to different processors
- For MPI to be useful, at least two or more blocks and at least three processors will be required.
- Often grids will arrive as multiple block grids. However, there are several reasons that additional block splitting will be required:
 - If the original mesh is not split into a sufficient number of blocks to efficiently use the processors available.
 - If the blocks are of disparate sizes, so that load balancing will be difficult.

Overview

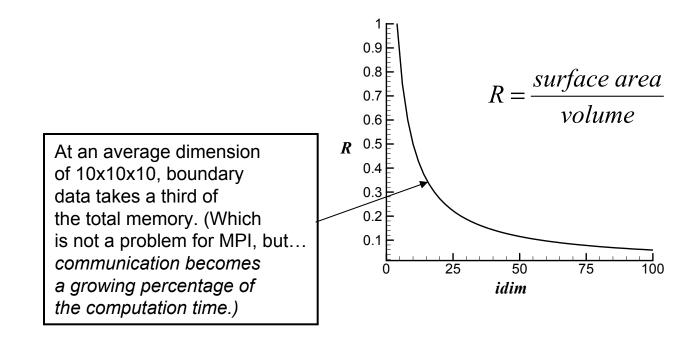


- Note, however, that there is a limit on the number of blocks for a given overall grid size for which efficient parallelization can take place.
 - Problem of growing communications between processors compared to processing per block (communication time).
 - Because CFL3D treats block boundaries explicitly, splitting into an ever increasing number of blocks amounts to making the code explicit. An increasing number of blocks means that an increasing number of sub-iterations will be required.
- The following illustrates the increasing communications with decreasing block sizes....



Problem of the humming bird versus the elephant

Consider the ratio of number of surface points to the total number of grid points as grid size diminishes. These results are based on a grid having equal idim, idim, kdim dimensions.



Overview

With the issues clearly in mind, there are times when splitting is useful...

- The tool 'splitter' is available with CFL3D for use in splitting blocks.
- It is created by performing the following command in the 'build' directory:

make splitter

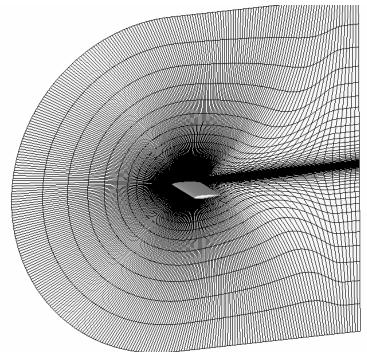
- The executable will be in the directory '~/cfl3dv6/build/split/seq/'.
- An example input can be found in the CFL3D version 6 web page.

Example: Splitting a single C-H grid



Lets consider again the BACT wing we have looked at previously. This grid has i,j,k dimensions 73 (spanwise) x 345 (streamwise) x 73 (normal to wing).

Suppose a 32 processor PC cluster is available for this problem. It would be useful to split this block into at least 24 blocks. However consideration must also be given to how many times each dimension can be split and still retain the proper dimensions to perform multi-grid computations.



Example: Splitting a single C-H grid



An acceptable block split can be obtained by requiring M, the number of split blocks, in the following computation

$$M = \frac{D-1}{d-1}$$

be an integer. D is the overall dimension of the un-split grid, and d is the proposed dimension of the split grid. For the current example, the j-dimension can be split with blocks having dimension of 9, 87 or 173.

$$M = \frac{345 - 1}{9 - 1} = 43$$
 , $M = \frac{345 - 1}{87 - 1} = 4$, $M = \frac{345 - 1}{173 - 1} = 2$

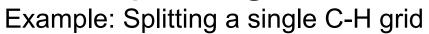
Example: Splitting a single C-H grid



Note that block dimensions of 87 or 173 will allow only 3 levels of multigrid, a dimension of 9 allows 4. We will chose a dimension of 87.

Similar computations for the idim = 73 and kdim = 73 lead us to chose 6 blocks in those directions with dimension of 13. This will result in a total of 144 blocks. This number of blocks will allow us to use 4, 24, 48 or 144 processors efficiently.

These computations result in 3 splits in the j-direction, 5 splits in the i-direction and 5 splits in the k-direction for a total of 13 splits. The input that performs these splits is shown in the next slide.





The splitter input file for this grid is

shown: INPUT (UNSPLIT) FILES cfl3d.inp ronnie.inp grid.unf

sd_grid.unf ICFLVER **IRONVER IGRDFMT ISDFMT OUTPUT (SPLIT) FILES**

cfl3d.inp_split ronnie.inp split grid_split.unf

sd_grid_split.unf ICFLVER **IRONVER IGRDFMT ISDFMT**

NSPLITS



Example: Splitting a single C-H grid

```
INPUT (UNSPLIT) FILES
cfl3d.inp
ronnie.inp
grid.unf
sd_grid.unf
                        IGRDFMT
ICFLVER
           IRONVER
                                     ISDFMT
       5
OUTPUT (SPLIT) FILES
cfl3d.inp split
ronnie.inp split
grid_split.unf
sd grid split.unf
ICFLVER IRONVER
                        IGRDFMT
                                     ISDFMT
```

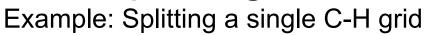
cfl3d.inp - cfl3d input file for the unsplit grid

ronnie.inp - ronnie input file for the unsplit grid, if not a patched case, enter the word **null**

grid.unf - grid file for the unsplit grid; can be formatted or unformatted

sd_grid.unf - sensitivity file for the unsplit grid NOTE: Currently not supported in Version 6; the same functionality is now handled via complex variables and a complex-valued grid file;

enter the word null





INPUT (UNSPLIT) FILES cfl3d.inp ronnie.inp grid.unf sd_grid.unf **ICFLVER** IRONVER **IGRDFMT ISDFMT** 5 **OUTPUT (SPLIT) FILES** cfl3d.inp split ronnie.inp split grid_split.unf sd grid split.unf ICFLVER IRONVER **IGRDFMT ISDFMT**

cfl3d.inp split

- cfl3d input file for the split grid

ronnie.inp split

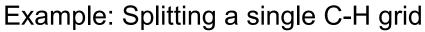
- ronnie input file for the split grid, if not a patched case, enter the word null

grid split.unf

- grid file for the split grid; can be formatted or unformatted

sd_grid_split.unf

- sensitivity file for the split grid NOTE: Currently not supported in Version 6; the same functionality is now handled via complex variables and a complex-valued grid file; enter the word **null**





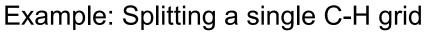
```
INPUT (UNSPLIT) FILES
cfl3d.inp
ronnie.inp
grid.unf
sd_grid.unf
ICFLVER IRONVER
                        IGRDFMT
                                    ISDFMT
      5
                  1
OUTPUT (SPLIT) FILES
cfl3d.inp split
ronnie.inp_split
grid_split.unf
sd grid split.unf
ICFLVER IRONVER
                       IGRDFMT
                                    ISDFMT
```

icflver

- = 4 the cfl3d input file is a version 4.1 type
- = -4 the cfl3d input file is a version 4.1hp type
- = 5 the cfl3d input file is a version 5/6 type

ironver

- = 0 ronnie input file is the old style, with all "from" blocks listed on one line
- = 1 ronnie input file is the new style, with each "from" block having it's own line NOTE: a value for ironver must always be entered, even if the case does not involve patched grids.





```
INPUT (UNSPLIT) FILES
cfl3d.inp
ronnie.inp
grid.unf
sd_grid.unf
ICFLVER IRONVER IGRDFMT ISDFMT
5 1 1 1 1
OUTPUT (SPLIT) FILES
cfl3d.inp_split
ronnie.inp_split
grid_split.unf
sd_grid_split.unf
ICFLVER IRONVER IGRDFMT ISDFMT
5 1 1 1
```

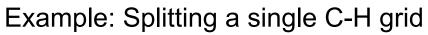
igrdfmt

- = 0 grid file is formatted
- = 1 grid file is unformatted

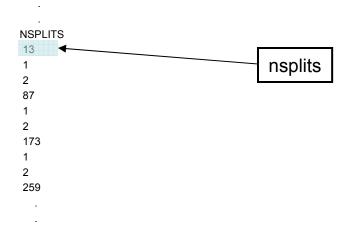
isdfmt

- = 0 sensitivity file is formatted
- = 1 sensitivity file is unformatted

NOTE: Currently not supported in Version 6; the same functionality is now handled via complex variables and a complex-valued grid file; however a value is still required - use 0 or 1

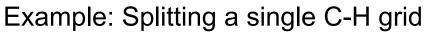




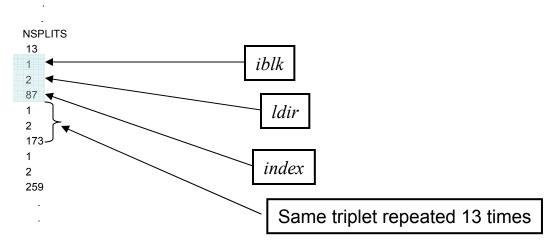


nsplits

- number of grid splits to perform (can be 0 in order to convert grid from formatted to unformatted or vice versa). Following the value of *nsplits*, *nsplits* triplets of integers must appear, one integer of the triplet per line....







iblk

- block number of the block to be split. NOTE: *iblk* always refers to the original, unsplit block number

ldir

- = 1 split in the *i*-direction
- = 2 split in the *j*-direction
- = 3 split in the *k*-direction

index

- split the block in the *ldir* direction at this value of the index





```
SPLITTER - CFL3D BLOCK AND INPUT FILE SPLITTER
    VERSION 6.X: Computational Fluids Lab, Mail Stop 128,
             NASA Langley Research Center, Hampton, VA
             Release Date:
                              MMM DD, YYYY.
memory allocation: 431.046108 Mbytes, double precision
input (unsplit) files
 cfl3d.inp
 null
 wbgrid.cfl
 null
icflver ironver igrdfmt isdfmt
   5
output (split) files
 cfl3d.inp_split
 null
 wbgrid_split.cfl
 null
icflver ironver igrdfmt isdfmt
```

Example: Splitter output



converting unsplit cfl3d input file to tlns3d map file checking dimensions... reading grid... grid: wbgrid.cfl block # 1: il= 73, jl= 345, kl= 73 number of splits = 13 split block coord index 87 173 259 13 25 37 6 49 61 Κ 13 Κ 25 Κ 37 11 Κ 49 12 1 1 Κ





```
49 61 259 345 49 61
30
           49 61 173 259 49 61
31
           49 61 87 173 49 61
                                                                          121
                                                                                      61 73 1 87 1 13
32
           49 61 1 87 49 61
                                                                          122
                                                                                      61 73 87 173 1 13
33
           37 49 1 87 49 61
                                                                          123
                                                                                      61 73 173 259 1 13
34
           37 49 87 173 49 61
                                                                          124
                                                                                      61 73 259 345 1 13
                                                                          125
                                                                                      49 61 259 345
           37 49 173 259 49 61
35
                                                                          126
                                                                                      49 61 173 259 1 13
           37 49 259 345 49 61
36
                                                                          127
                                                                                      49 61 87 173 1 13
37
           25 37 259 345 49 61
                                                                          128
                                                                                      49 61
                                                                                             1 87 1 13
38
           25 37 173 259 49 61
                                                                          129
                                                                                      37 49 1 87 1 13
39
           25 37 87 173 49 61
                                                                          130
                                                                                      37 49 87 173 1 13
                                                                          131
                                                                                      37 49 173 259 1 13
40
           25 37 1 87 49 61
                                                                          132
                                                                                      37 49 259 345 1 13
41
           13 25 1 87 49 61
                                                                          133
                                                                                      25 37 259 345 1 13
42
           13 25 87 173 49 61
                                                                          134
                                                                                      25 37 173 259 1 13
43
           13 25 173 259 49 61
                                                                                      25 37 87 173 1 13
                                                                          135
44
           13 25 259 345 49 61
                                                                          136
                                                                                      25 37 1 87 1 13
                                                                          137
                                                                                      13 25
                                                                                             1 87 1 13
45
            1 13 259 345 49 61
                                                                                      13 25 87 173 1 13
                                                                          138
46
            1 13 173 259 49 61
                                                                          139
                                                                                      13 25 173 259 1 13
47
            1 13 87 173 49 61
                                                                          140
                                                                                      13 25 259 345 1 13
            1 13 1 87 49 61
                                                                          141
                                                                                       1 13 259 345 1 13
                                                                          142
                                                                                       1 13 173 259 1 13
                                                                          143
                                                                                       1 13 87 173 1 13
                                                                                             1 87 1 13
                                                                                       1 13
```

split-grid basic dimensions are multigridable to ncg = 1

Input points: 1838505 Ouput points: 2117232

Block Splitting and MPI



Notes regarding use:

- IF A LIMITER IS DESIRED, USE IFLIM=4. This will allow for consistent results with block splitting; iflim=3 is not recommended iflim=4 is basically a correct implementation of iflim=3 for multiple blocks, and should now be viewed as the recommended limiter for any case that needs one.
- Also, for exact consistency between split and unsplit grids, version 5 emulation (i.e. "Install -v5) should not be used. Version 5 (and earlier versions) made an approximation for cell volumes at 1-1 block interfaces that has been eliminated in version 6 in favor of the exact treatment.
- The input file part of the splitter works by first converting the unsplit CFL3D input file to a TLNS3D map file, splitting the TLNS3D map file, then converting the split TLNS3D map file back to a CFL3D input file.

Block Splitting and MPI



Notes (...continued):

- Caveats: The conversions from the CFL3D input file to a TLNS3D map file are not perfect! The user is urged check the resulting split CFL3D input (and patch) files.
 - A useful check before actually splitting the files is to run this splitter with the number of splittings = 0, and the output grid file as null. Running splitter in this way will cause to code to go through the translations, but the "split" files will have the same numbers of blocks, and the "split" grid will not be output.
 - A "diff" or "gdiff" will point to translation-induced differences that should be easier to sort out than when coupled with true splitting. Note that the 2-step process almost always results in a *reordering* of some boundary condition segments.



MPI requires one processor for overhead. For example if a 32 processor cluster is employed, and there are 28 blocks to be computed on 28 processors, then the command line will read:

```
mpirun –np 29 cfl3d_mpi < cfl3d.inp &
```

 You may want to verify the correct procedure for running mpi code on your platform (e.g. some mpp's use -n instead of -np)



- Because version 6 has dynamic memory allocation, there is no *requirement* to run precfl3d before you can run cfl3d. However, you may still find it useful to do so in order to assess how much memory will be required to run the case at hand, allowing you to determine whether a particular problem can fit within the memory of the machine, or to determine the appropriate queue in which to submit the job.
- The usage of precfl3d has changed slightly from previous versions: you must now specify the number of processors in addition to the input file, for example:

precfl3d -np num_procs < cfl3d.inp &</pre>

where **num_procs** is the total number of processors, including the host. When running on a single processor, that processor is the host, so num_procs=1 will suffice to assess the memory requirements for the sequential version of the code.

An important reason why you may want to run precfl3d before running the parallel version of the code is that for num_procs > 1, precfl3d will output an auxiliary file called ideal_speedup.dat.
 This file will list the best possible speedup you could hope to achieve for the current case, using various numbers of compute processors, ranging from 1 to the number of zones in your grid.



The BACT case with 144 blocks was run on 24 processors (-np 25). In the 'precfl3d.out' file the following information is contained:

BLOCK TO NODE MAPPING no. of blocks = 288 no. of nodes = 24 block node



```
SUMMARY OF STORAGE REQUIREMENTS - W + WK ARRAYS
sequential version:
     permanent array w requires 131825665 (words)
     temporary array wk requires 2681342 (words)
     temporary array iwk requires 187820 (words)
parallel version, per node:
     permanent array w requires 5506908 (words)
     temporary array wk requires 1500235 (words)
     temporary array iwk requires 187820 (words)
>>> Estimate for mwork
                       (seguential) = 134507007 <<<
>>> Estimate for mworki
                      (sequential) = 187820 <<<
>>> Estimate for mwork (per node, parallel) = 7007143 <<<
>>> Estimate for mworki (per node, parallel) = 187820 <<<
>>> Parallel code sized for 24 nodes, min. (+host)
****************
```



In the 'cfl3d.out' file the same information is found plus the following contained at the end:

computational rate by mesh sequence (based on wall time): iseq 1 181.13 microseconds/cell/time step 90.56 microseconds/cell/subiteration

timing for complete run - time in seconds

node	user	system	total	wall clock
0	10.15	17.60	27.75	325.00
1	3.64	0.55	4.19	228.00
2	5.37	0.92	6.29	325.00
3	3.90	0.52	4.42	228.00
4	5.36	0.87	6.23	325.00
5	5.85	1.14	6.99	324.00
6	4.54	0.89	5.43	228.00
7	4.38	0.83	5.21	227.00
8	4.03	0.79	4.82	226.00
9	4.31	0.70	5.01	228.00
10	6.08	1.00	7.08	325.00

```
5.17
                          227.00
     4.40
            0.77
12
    4.19
            0.65
                          227.00
                   4.84
13
     4.20
            0.74
                   4.94
                          226.00
     4.42
                          225.00
14
            0.66
                   5.08
15
     4.25
            0.81
                   5.06
                          226.00
     4.35
            0.68
                   5.03
                          225.00
17
     4.08
            0.83
                   4.91
                          225.00
18
     4.22
            0.87
                   5.09
                          225.00
     4.35
                          225.00
19
            0.66
                   5.01
20
     4.17
            0.66
                   4.83
                          225.00
     3.78
            0.55
                   4.33
                          224.00
     3.59
            0.49
                   4.08
                          225.00
     3.58
                          224.00
            0.51
                   4.09
     3.40
            0.40
                   3.80
                          224.00
```

total: 114.59 35.09 149.68

total run (wall) time = 0 hours 3 minutes 44 seconds

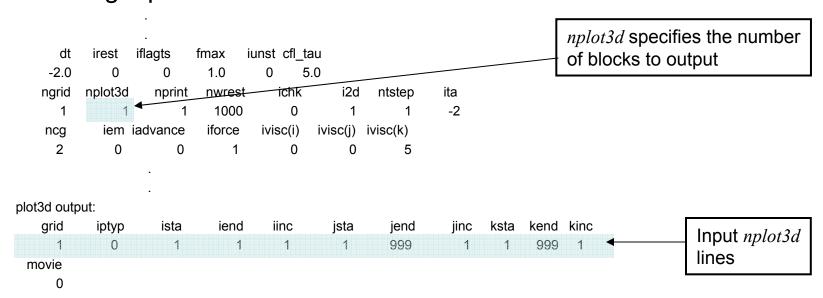
memory for cfl3d has been deallocated

Flow Field Visualization



Plot3D output

CFL3D is capable of creating Plot3D files of the grid and flow field. Specification of the region of the flow field for output is found in the following input lines:



If nplot3d < 0, then the Plot3D files are automatically set to include all solid Surfaces (no field points) for 3D cases or all field points for 2D cases

Flow Field Visualization



Plot3D output

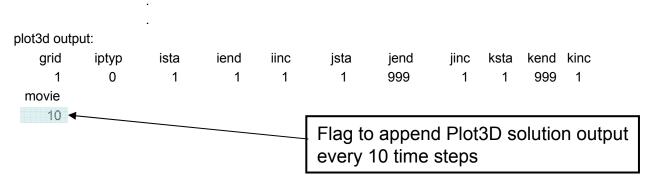
		•								
plot3d outp	ut:	•								
grid	iptyp	ista	iend	iinc	jsta	jend	•	ksta		nd kinc
1 movie	0	1	1	1	1	999	1	1	999	99 1
0										
Grid	-	- Desi	gnated	d grid	numb	er to b	e ou	tput		
iptyp	:	= 0 -	grid p	ooint t	уре –	grid fil	le an	d Q	file	e output
	= 1 - cell center type – grid file and Q file output									
	= 2 - cell center type - grid file and turbulence file output (ivisc > 1 only)									
	;	> 2 -	near	est vis	scous	•	r dire			ction file output ($iptype = 3$ – minimum distance to istance ($ivisc > 1$ only), $iptype = 4$ – eddy
ista, jst	a, ksta	<i>a</i> -	startin	g indi	ces in	the i,j	k dir	ecti	ons	3
iend,jei	nd,ken		-	_		Ū			•	(note that if these values are set higher than them to the block dimensions)
iinc,jin	c,kinc	-	increr	nent i	n the i	<i>,j,k</i> dir	ectio	ns		

Note: Setting ista = iend = iinc = 0, etc... is a short hand way of specifying the entire range.

Flow Field Visualization



Movie output



Note that one gird file and one solutions file are generated.

Movie = 0 no output of intermediate solutions (if nplot3d > 0), then a single solution is written at the end of the run.

Movie > 0 output of additional solutions every *movie* iterations (time steps)

Movie < 0 output of the initial flow field at the beginning of the run and output of additional solutions every *movie* iterations (time steps)

Caution: Use with care. Plot3D file will get very large very quickly.

The tool 'moovmaker' will read the plot3D solution and grid file and create a movie for a 2D flow field in which the 3rd dimension will be time. The grid output by 'moovmaker will be called 'g.bin' and the solution file will be called 'q.bin'. Creating these files will allow animating the 3rd dimension (time) to produce a movie of the flow field.



- Get_FD.F
 - This program reads two CFL3D restart files and calculate finite differences of force and moment coefficients; it is used to validate complex-variable approach for determining solution derivatives.
- INGRID_to_p3d.F
 - This program converts PEGSUS 4.x INGRID file to a PLOT3D file that can be used in CFL3D. Note that the INGRID file must correspond to grid points rather than "augmented" cell centers.
- XINTOUT_to_ovrlp.F
 - This program converts the XINTOUT overset grid interpolation file from PEGSUS to the ovrlp.bin file used by CFL3D.
- cfl3d_to_pegbc.F
 - This program creates a peg.bc.raw file for use with PEGSUS 5.x.
- cgns_to_cfl3dinput.F
 - This program reads a CGNS file and creates a PLOT3D-type grid as well as a best-guess for a CFL3D input file.



everyother_xyz.F

 This program reads a grid and creates an every-other-point grid. This can be useful in combination with the program v6inpdoubhalf.F, in order to reduce the required CFL3D runtime memory when you are only running on a coarser-level grid (and not taking it up to the finer level(s).

grid_perturb.F

This program generates a real-valued grid (PLOT3D multiblock form) by reading in a real-valued grid (PLOT3D multiblock form) and a corresponding real-valued matrix of grid-sensitivity derivatives (PLOT3D multiblock function file form, with 3*ndv variables for the x,y,z components of the ndv design variables). The code Get_FD.F may be used with the two restart files to determine d(Cl)/d(DV), d(Cd)/d(DV), etc.

grid_perturb_cmplx.F

This program generates a complex-valued grid (PLOT3D multiblock form) by reading in a real-valued grid (PLOT3D multiblock form) and a corresponding real-valued matrix of grid-sensitivity derivatives (PLOT3D multiblock function file form, with 3*ndv variables for the x,y,z components of the ndv design variables). The output grid may be read into the complex version of CFL3D (cfl3dcmplx_mpi or cfl3dcmplx_seq) to determine the solution derivatives with respect to the chosen design variable.



initialize_field.F

 This program creates a restart.bin restart file in which you can specify specific initial conditions, region by region. This can be useful when "freestream everywhere" is not a desirable initial condition.

moovmaker.F

 This program reads the PLOT3D files output by CFL3D when the MOVIE parameter is used for 2-D datasets (or 3-D datasets surface-only), and creates new PLOT3D files with time as the third (k) direction.

p3d_to_INGRID.F

 This program converts either PLOT3D or CFL3D type grids into either INGRID type grids that can be used with PEGSUS 4.x, or PLOT3D type grids that can be used with PEGSUS 5.x.
 The converted grids can contain either the grid points as given in the input grids, or "augmented" cell centers of the input grids.

p3d to cfl3drst.F

This program reads PLOT3D files and creates an approximate restart.bin restart file. This
can be useful if: (1) you are given a PLOT3D Q-file from another code, and you wish to use it
as a basis for starting CFL3D, or (2) you have lost the CFL3D restart file, but you still have
the PLOT3D Q-file.



- plot3dg_to_cgns.F
 - This program reads a PLOT3D grid file and a CFL3D input file and creates a CGNS file (with grid, BC, and 1-to-1 connectivity information in it).
- v6_restart_mod.F
 - This program reads a restart.bin restart file and manipulates it. It can switch between unformatted and formatted (which is useful if you need to transfer the restart file to a machine of different architecture). It can also write out the restart file either the same size, half the size, or double the size. Going to half size is useful if one wishes to restart from a fine grid solution and run on a coarser level. User can choose to coarsen/refine only particular index directions, if desired. The program cannot both coarsen and refine different directions simultaneously.
- v6inpdoubhalf.F
 - This program reads a CFL3D input file and creates a new input file appropriate for a grid of either half or double the size. This can be useful in combination with the program everyother_xyz.F when running on coarser grid levels, and you wish to reduce the run-time memory required.

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Summary



- CFL3D is a general purpose production-level CFD code for fluid dynamics, with many capabilities and options.
- This tutorial has summarized many of the newest features of the code, and also has explained in detail how to set up and run it for general cases.
- Particular focus has been given to CFL3D's upgraded deforming mesh and aeroelastic analysis capabilities.

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14. ABSTRACT

This document contains the course notes on the computational fluid dynamics code CFL3D version 6.4. It is intended to provide from basic to advanced users the information necessary to successfully use the code for a broad range of cases. Much of the course covers capability that has been a part of previous versions of the code, with material compiled from a CFL3D v5.0 manual and from the CFL3D v6 web site prior to the current release. This part of the material is presented to users of the code not familiar with computational fluid dynamics. There is new capability in CFL3D version 6.4 presented here that has not previously been published. There are also outdated features no longer used or recommended in recent releases of the code. The information offered here supersedes earlier manuals and updates outdated usage. Where current usage supersedes older versions, notation of that is made. These course notes also provides hints for usage, code installation and examples not found elsewhere.

15. SUBJECT TERMS

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